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(71) Applicant (for all designated States except US): UNITED STATES OF AMERICA, represented by THE SECRE-TARY OF THE ARMY [US/US]; Intellectuel Property Counsel of the Army, Office of The Judge Advocate General, DA, Suite 400, 901 North Stuart Street, Arlington, VA 22203-1837 (US).

(72) Inventors; and

- (75) Inventors/Applicants (for US only): REID, Robert, H. [US/US]; 10807 McComas Court, Kensington, MD 20895 (US). SADEGH-NASSERI, Scheherazade [US/US]; 13600 Straw-Bale Lane, Damestown, MD 20878 (US). WOLFF, Marcia [US/US]; 9850 Hollow Glen Place, Silver Spring, MD 20910 (US). NAUSS, Jeffrey, L. [US/US]; 142 Martha Lane, Fairfield, OH 45014 (US).
- (74) Agent: NORRIS, Jerome, J.; Suite 1250, 1401 New York Avenue, N.W., Washington, DC 20005 (US).

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(54) Title: MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

(57) Abstract

Assay methods for determining whether a peptide is likely to be immunogenic are based on a computer modeling of binding to a Class II MHC DR1 receptor. This is confirmed by competitive inhibition binding assays. The peptides are useful for eliciting an immune response for vaccination or the production of antibodies or T-cells.

Applicants: Alexander Gad and Dora Lis

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Exhibit 11

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MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

Government Interest

The invention described herein may be manufactured, licensed and used by or for governmental purposes without the payment of any royalties to us thereon.

Cross Reference

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This application is a continuation-in-part of U.S. Patent application Serial No. 08/064,559, filed May 21, 1993, and the present application incorporates U.S. Patent Application Serial No. 08/064,559 in its entirety by reference.

15 Field of the Invention:

This invention relates to a means of predicting potential of a peptide for eliciting immune response.

Background of the Invention:

Among the numerous steps required for an immunological response to occur is the presentation of the antigen by macrophages to the B-cell or T-cell. This presentation is mediated by the Class I and Class II major histocompatibility complex (MHC) molecules on the surface of the cell. The MHC molecules hold antigens in the form of the peptide fragments and together with the receptor molecule on the T-cells, form a macromolecular complex that induces a response in the T-cell. Therefore, a necessary step in an immune response is the binding of the antigen to the MHC.

Recent single crystal X-ray structures of human and murine Class I MHC's have been reported. Analysis of these crystal structures have shown that antigenic peptides lie in the so-called binding cleft for presentation to the T-cell. This cleft is formed by α_1 and α_2 domains and by β -strands from each domain forming the floor. Furthermore, the sequence polymorphism among Class I molecules can result in alterations of the surface of the cleft forming different pockets. Peptide side chains may insert into these pockets. Thus, different pockets may interact with different side chains. This implies the mechanism for the peptide specificity of Class I MHC's. Peptides bound to the Class I MHC's in the crystal structures were found to have both the amino and carboxy termini tightly held by the MHC. There were few interactions near the middle of the cleft. Hence the bound peptide is allowed to bend slightly in the center. The observed binding mode helped to explain the apparent partial specificity of peptide sequence and the allowed variation in peptide length found among peptides isolated from Class I MHC's.

The precise mode of binding of peptides to Class II MHC molecules is less clear. While a single crystal X-ray diffraction structure for the HLA-DR1 MHC has been shown, the coordinates have remained unavailable. However, currently available theoretical and experimental results help form a hypothesis that the binding of a peptide to Class II MHC is similar to that observed with Class I. First, it is noted that the Class II binding cleft is structurally similar to

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that of Class I. This was concluded based upon a sequence analysis of 26 Class I and 54 Class II amino acid sequences.

Unlike with Class I molecules, self-peptides isolated from murine I-Ab and I-Eb, from murine I-Ad and from human HLA-DR1 molecules were found to be varied in size (13 to 25 residues long). The peptides isolated from the murine I-Ab and I-Eb molecules had heterogenous carboxy termini while those from I-Ad and HLA-DR1 had ragged termini at both ends. The varying lengths indicate that the amino and carboxy termini of the peptides were not critical for the binding. One or both termini may protrude from the binding site and be available for further processing. The residues critical for binding were proposed to be at the ends of the peptide as opposed to the center.

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Summary of the Invention:

It is the purpose of this invention to provide a method for preliminary screening of peptides for ability to elicit an immune response. Structural homology techniques were used to model a receptor (the Class II MHC is exemplified). This model makes it possible to preliminarily screen peptides for antigenic properties. By modifying the peptide to "fit" into the receptor it is possible to identify methods of rendering non-immunogenic peptides immunogenic.

The preliminary screening of peptides for immunogenicity comprises the steps of (1) creating a molecular model of a receptor followed by minimizing the model created, 2) modeling a peptide to

be tested and minimizing the model of the peptide, then testing the

fit of the model of the peptide into the model of the receptor to produce a composite minimized receptor/minimized peptide model.

Upon finding an acceptable fit, the peptide may then be screened by a binding assay for actual binding to Class II MHC as a further test for immunogenicity.

It has been found that when the model of the peptide can not be fitted into the model of the receptor, the peptide will lack immunogenicity. While not all peptide models which can be made to "fit" into to model of the receptor will be effective as immunogens, the screening methods of the invention may make it possible to avoid undue biological testing of inappropriate peptides. By using the model, it is also possible to alter peptides to accommodate the receptor. Hence, the invention has both predictive and drug design applications.

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Brief Description of the Figures:

Fig. 1 shows the HLA-aw68 α_1 and α_2 domains with DR1 α_1 and β_1 domains.

Figs. 2-30 are a printout of the minimized coordinates of the 20 receptor.

Figs. 31 and 32 shows the effects of various peptides inhibiting the binding of labeled hemagglutinin in a competitive binding assay.

25 <u>Detailed Description of the Invention:</u>

In order to understand and better predict peptide interaction with Class II MHC's and as an aid for synthetic peptide vaccine design, a structural homology model of HLA-DR1 molecule was made

using the Class I HLA-aw68 as a reference molecule. For purposes of this analysis, numerous conserved residues were aligned leading to a proposed three-dimensional model for the Class II structure very similar to that of Class I. This model retained the overall conformation of a Class I MHC and agreed with a considerable amount of the published data. Furthermore, peptides shown to bind to DR1 were docked in the binding cleft of the model and analyzed. The results agree with the experimental binding data presented here. Hence, it is shown that the structural homology model reported here is useful for screening Class II MHC functionality.

It had been hypothesized that few peptide residues may be required for binding to DR1. By substituting residues into the influenza hemagglutinin 307-319 T-cell epitope (HA) it had been determined that a single tyrosine at 308 was required for binding. A synthetic peptide with the tyrosine at position 308 and a lysine at 315 was found to bind DR1 as well as the native peptide. Hence, it was concluded that few peptide residues determine the high affinity binding to DR1.

The peptides produced according to the present invention may be used alone or chemically bound to another peptide and/or carrier in order to elicit an immune response. An immune response is elicited by administering a peptide to an animal in an effective dose and by an effective route of administration. Typically the peptide will be administered with an immunologically acceptable carrier. The routes of administration, dosages, times between multiple administrations will be based on the particular peptide and are standard operations of those skilled in the art.

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Of particular interest are peptides from pathogenic microorganisms and neoplasms. In such an example, a vaccine may be formed with the peptide and any known immunological carrier and may be administered prophylactically or therapeutically. The immune response may be elicited for a number of reasons other than for prophylaxis or therapy such as increasing antibody production for the harvesting of antibodies, or increasing specific B-cell or T-cell concentration for the production of hybridomas or cellular therapy.

The choice of host animals is limited only to those capable of an immune response. Preferred hosts are mammals, more preferred are humans.

The vaccine may contain plural peptides with each peptide corresponding to the same or different antigens. The peptides may be used unbound or they may be chemically bound to another peptide or an unrelated protein or other molecule. A preferred vaccine preparation contains a plurality of peptides chemically bound to a larger more immunogenic peptide.

The peptide may be adsorbed, bound or encapsulated in a biodegradeable microsphere, microcapsule, larger carrier or a combination of these. The carrier may have a slow or controlled release property thereby releasing the peptide under appropriate conditions and times for enhanced immunization. This is particularly important when administering the peptide orally where stomach acid can degrade the peptide.

Another embodiment of the present invention is to modify the amino acid sequence of a peptide to enhance its immunogenicity.

This is done by modifying the natural peptide sequence to bind to

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the Class II MHC receptor DR1 with superior binding affinity for a Class II MHC receptor DR1 than the natural peptide sequence. This modified peptide is considered a synthetic peptide. Alternatively, the sequence may be modified to have a greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1.

Many amino acid changes are acceptable in the formation of a synthetic peptide. The changes may be for similar types of amino acids such as leucine for isoleucine or they may be for diverse types such as tyrosine for lysine.

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Materials and Methods:

The structural homology model for the DR1 Class II MHC was constructed using the QUANTA molecular modeling package (vision 3.2, Molecular Simulations, Inc., Burlington, MA) with the CHARMM and Protein Design modules. After alignment of the sequences as described below, gaps and loops were energy minimized using 100 steps of steepest descents minimization followed by 100 steps of adopted basis set Newton-Rapheson (ABNR) minimization. were closed using a fragment database from a selected set of high-resolution crystal structures. The resulting structure was minimized in vacuo using 1000 steps of steepest descents followed by an additional 1000 steps of ABNR minimization. A distance related electrostatic function was used in all calculations with a dielectric constant of 1.0. Non-bound parameter lists were updated every 20 steps with a cutoff distance of 15.0Å. Non-bonded calculations were performed using a shifted potential function between 11.0Å and 14.0Å. An extended atom set was used with only

polar hydrogen atoms specifically placed. There were no explicit hydrogen bond energy calculations performed.

All peptides were initially modeled using QUANTA in an extended chain conformation and subjected to 500 steps of ABNR minimization. The resulting structures remained essentially in extended chain conformations. Individual peptides were manually docked in several different orientations into the binding cleft region of the minimized DR1 structure. The resulting bimolecular complex was subjected to 5000 steps of steepest descents minimization with non-bonded interactions updated every five steps. After minimization, bound peptides remained essentially in extended chain conformations. The lowest energy complexes for each peptide were selected for further analysis.

The selected peptide and DR1 complexes and the minimized DR1 model were subjected to the following molecular dynamics regimen: 300 steps of heating to 300°K, 600 steps of equilibration at 300°K, and 1100 steps of production dynamics. During this simulation, the DR1 C α atoms were constrained in their starting positions. All non-bonded interaction parameters were as stated for the minimization procedure. The lowest energy structure during the course of the production dynamics was selected and subjected to the 5000 step minimization procedure described previously with the C α restraints removed. The resulting structures were used for the binding energy calculations and for hydrogen bonding analysis.

Hydrogen bonds were determined using the QUANTA default parameters. Maximum allowed distances were 2.5Å between a hydrogen and the acceptor atom and 3.3Å between the donor and acceptor atoms.

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The minimum angle allowed between any set of atoms forming a hydrogen bond was 90°.

Competitive Inhibition Binding Assay:

Was labeled with ¹²⁵I. The labeled HA peptides were then allowed to interact with purified DR1 molecules during incubation to allow formation of peptide/DR1 complexes. After incubation, the peptide/DR1 composition was exposed to a native gel for chromatographic separation or passed through a spun column to separate labeled peptide/DR1 complex and free labelled peptide. When unlabeled peptides were added before incubation of labeled HA peptides and DR1, and if the unlabelled peptides had capacity for binding to DR1 simultaneous with ¹²⁵I-HA, there was a resultant decrease in radioactive signal associated with the DR1. The extent of this decrease directly related to the binding capacity of the unlabeled unknown peptide.

Structural Homology Model for the DR1 Molecule:

The structural homology model was created, the reference molecule being the crystal structure of HLA-aw68. The HLA-aw68 coordinates and subsequent sequence were obtained from the entry 2HLA in the Brookhaven Protein Data Bank released January 15, 1991, which is incorporated herein by reference. The sequence for the DR1 molecule was for the α_1 domain was reported by Klein and for the β_1 domain, the study reported by Todd et al. (Nature 329, 599 (1987)).

The sequence alignment is based on Brown et al. (Nature 332, 845 (1988)). The complete alignment and numbering scheme for both

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are seen in Figure 1. The Class II, β_1 and Class I α_2 domains regions were conserved with some variations at the ends where the two MHC's have different loop regions. The fourth B-strand in the α_1 domain of HLA-aw68 (residues 30-38) is disrupted in the DR1 model. Only three residues are in a β -sheet conformation, probably due to the inserted glycine at position 28 before the strand and the large deletion in the loop region immediately after the strand. two alpha-helical regions are clearly maintained. Both helices have been observed to be discontinuous in the Class I molecules and are similar in the DR1 model. The α_1 domain helix is long and curves from residues 49α to 76α without significant disruption. essentially a single continuous helix. However, the α_2 helical region is broken into two separate helices as with the Class I molecules. A short helix (52-63) is separated from a longer helix (68-94) by a deformed region without secondary structure. This deformation is more pronounced in the DR1 model as opposed to the Class I molecules due to an insertion.

Influenza Hemagglutinin Peptide with DR1:

The amino acid residues 307-319 of influenza hemagglutinin (Pro-Lys- Tyr-Val-Lys-Gln-Asn-Thr-Leu-Lys-Leu-Ala-Thr) make up a well-documented linear T-cell epitope which has been shown to be HLA-DR1 restricted. With the demonstration that the influenza hemagglutinin epitope (referred to as the HA peptide) binds DR1, it was chosen to be modeled into the binding cleft.

The peptide was initially inserted into the cleft so that Leu 11 HA was in the vicinity of the hydrophobic pocket. This allowed Asn 7 to be near the middle charged and polar groups of the cleft.

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The remaining residue of the motif (Lys 2) was near the vicinity of the remaining charged and polar residues at the end of the cleft.

The only adjustment to the starting conformation was a slight rearrangement of the terminal peptide proline and Tyr 3 to alleviate obvious bad contacts.

After the energy minimization of the bimolecular complex, the total energy was reduced to 483 kcal/mol. This reduction in energy was accomplished by alleviation of several bad contacts and also be formation of several hydrogen bonds. The sticking feature of this mode is lack of hydrogen bonds in the carboxy terminal half of the peptide. Only one hydrogen bond is identified between the backbone carbonyl group of Leu 9 and the side chain of the β_1 Asn 77. In contrast, the amino terminal half has eleven identified interactions. Four of these interaction involve the peptide backbone residues Tyr 3, Val 4, and Gln 6. The remainder involve the side chains of Lys 2, Tyr 3, Lys 5 and Gln 6. Interestingly, Lys 5 is involved in more interactions (three) than Lys 2 (only 2). No interactions were observed as anticipated with Asn 7. Instead, it was the glutamine at position 6 donating a hydrogen bond to the α_1 Asn 62. No interactions were observed for the amino and carboxy termini.

HA-YK Peptide with DR1:

The binding of the HA-YK peptide (Ala-Ala-Tyr-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Lys-Ala-Ala) to the DR1 model was tested. In aligning the peptide in the cleft, it was deemed logical to insert the tyrosine residue into the hydrophobic region of the binding cleft. The lysine would then be in position to interact with the

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hydrophilic groups in the other half of the cleft. The resulting peptide orientation is the opposite that used for the HA and the CS3 (defined below) peptides. With the peptide oriented as described, the final docking position for the peptide was unclear. The hydrophobic pocket is quite large, and, at least in this model, could accommodate the peptide tyrosine in a number of positions by sliding the peptide lengthwise through the cleft. However, repositioning the peptide also repositions the lysine. There were primarily two positions for the lysine: one with the lysine inside the cleft and the second with it outside. Of the two positions, the former was the lower in energy by 46 kcal/mol and had the greater number of interactions with the protein (11 vs. 7). Thus, the preferred orientation of the peptide appears to be with the lysine inside the binding cleft region.

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CS3 subunit Pilin Peptide with DR1:

The suspected T-cell epitope for CS3 pilus subunit 63-78

(Ser-Lys-Asn-Gly-Thr-Val-Thr-Trp-Ala-His-Glu-Thr-Asn-Asn-Ser-Ala)

was modeled with the DR1 molecule. The peptide was inserted with

lysine inside the cleft in the hydrophilic region. This placed the

Thr 5 in the center of the binding cleft and the tryptophane

(residue 8) near the hydrophobic region. The resulting minimized

model had ten interactions between the peptide and the protein,

three interactions with the peptide backbone and five with the

peptide side chains. The remaining two were with the amino terminal

of the peptide. All of the interactions were in either the first

three residues, His 10 or Glu 11 in the peptide. No interactions

were observed in the center of the cleft or residues four through nine.

CFA/1 with DR1:

A peptide identified as CFA/1 (colonization factor antigen)

(Val-Gly-Lys-Asn-Ile-Thr-Val-Thr-Ala-Ser-Val-Asp-Pro) was prepared

and an attempt was made to "fit" the molecule into the cleft of the

DR1. The lysine at position 3 prevented insertion of the peptide.

10 Results:

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The peptides chosen to dock in the DR1 model are shown in Table

1. The peptides were docked manually in several orientations into
the DR1 model. The peptides were then tested in biological binding
assays with the following results:

Table I

Peptide	Molecular Model predicted binding	Binding in the bioassay	
HA (influenza hemagglutinin)	Yes	Yes	
HA-YK (synthetic peptide)	Yes	Yes	
CS3 Pilin subunit	Yes	Yes	
CFA/1	No	No	

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Quantitative measurement of the inhibition of CS3 63-78 and HA 306-318 as compared to controls is shown in Fig. 31.

The binding energy was calculated as the difference between the final DR1 and peptide complex and the sum of the energies for the minimized DR and peptide models individually. The data is shown in Table II.

Table II.

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Peptide	Protein	Residues	Sequence	Binding Energy (kcal/mol)
НА	Influenza hemagglutinin	306-318	PKYVKQNTLKLAT	-283
HA-YK	synthetic peptide		ААҮАААААКАА	-216
CS3	CS3 pilin subunit	63-78	SKNGTVTWAHETNNSA	-245

CS6α and CS6ß with DR1

Colonization factor antigen IV (CFA/IV is an antigen on the surface of many enterotoxigenic *E. coli* one component of which is CS6. CS6 has two major subunits and a number of minor subunits. Several peptides from CS6 have been sequenced and assayed for potential inhibition of radiolabeled HA (306-318)/DR1 complex as a measure of immunogenicity. The sequences of the subunits are shown in Table III.

Table III.

Peptide	Amino Acid Residues	Sequence		
CS6α6	63-75	DEYGLGRLVNTAD		
CS6α7	80-92	IIYQIVDEKGKKK		
CS6α8	111-123	LNYTSGEKKISPG		
CS6ß1	3-15	WQYKSLDVNVNIE		
CS6ß2	42-54	QLYTVEMTIPAGV		
CS6ß3	112-124	TSYTFSAIYTGGE		
CS6ß4	123-135	GEYPNSGYSSGTY		
CS6ß5	133-145	GTYAGHLTVSFYS		

These peptides were assayed for inhibition of radioactively labeled HA(306-318)/DR1. The results are demonstrated in Fig. 32.

The foregoing description of the specific embodiments reveal the general nature of the invention so that others can, by applying current knowledge, readily modify and/or adapt for various applications such specific embodiments without departing from the generic concept, and, therefore, such adaptations and modifications should and are intended to be comprehended within the meaning and range of equivalents of the disclosed embodiments. It is to be understood that the phraseology or terminology employed herein is for the purpose of description and not of limitation.

All references mentioned in this application are incorporated by reference.

We Claim:

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1. A method of preliminarily screening peptides for immunogenicity comprising the steps of:

- 1) creating a molecular model of receptor DR1 Class II MHC and minimizing the model of the DR1;
 - 2) modeling a peptide to be tested and minimizing the model of the peptide; and
- 3) testing fit of model obtained in step 2 into the model obtained in step 1 to produce a composite receptor/peptide model.
 - 2. A computerized model comprising a model of the DR1 molecule having fitted in a cleft therein a model of a peptide.
- 3. A method of claim 1 wherein, additionally, the receptor/peptide model is subjected to computer-simulated heating.
 - 4. A method of claim 1 further comprising, assaying the peptide by competitive inhibition binding to a Class II MHC receptor DR1.
 - 5. A minimized peptide capable of binding to a Class II MHC receptor DR1 and inhibiting the binding of HA (306-318).
- 6. A synthetic peptide, wherein the amino acid sequence of the
 minimized peptide according to claim 5 has been modified to have a
 superior binding affinity for a Class II MHC receptor DR1 to form at
 least a portion of the synthetic peptide.

7. A synthetic peptide, wherein the amino acid sequence of the minimized peptide according to claim 5, has been modified to have greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1 to form at least a portion of the synthetic peptide.

- 8. A synthetic peptide according to claim 6, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 9. A synthetic peptide according to claim 7, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 10. A synthetic peptide according to claim 8, wherein said uncharged amino acid is alanine.
 - 11. A synthetic peptide according to claim 9, wherein said uncharged amino acid is alanine.
- 12. A minimized peptide according to claim 5, wherein the sequence is selected from the group consisting of PKYVKQNTLKLAT, AAYAAAAAKAA and SKNGTVTWAHETNNSA.
- 13. A minimized peptide according to claim 5, wherein the sequence is contained in a CFA.

14. A minimized peptide according to claim 13, wherein the sequence is selected from the group consisting of DEYGLGRLVNTAD, IIYQIVDEKGKKK, LNYTSGEKKISPG, WQYKSLDVNVNIE, QLYTVEMTIPAGV, TSYTFSAIYTGGE, GEYPNSGYSSGTY and GTYAGHLTVSFYS.

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- 15. A vaccine comprising:
 - a minimized peptide according to claim 5; and an immunologically acceptable carrier.
- 10 16. A vaccine comprising:
 - a synthetic peptide according to claim 6; and an immunologically acceptable carrier.
 - 17. A vaccine comprising:
- a synthetic peptide according to claim 7; and an immunologically acceptable carrier.
 - 18. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 15.
 - 19. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 16.

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20. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 17.

av68 DR1 #,	CSUSHRYF 9 YISVSRPGRG 19 EPRPIAVGYV 29 D DTGFVRPD 38 SDAASQRHEP 48 RAPHIEQUOR 1 CSUSHRYF 9 YISVSRPGRG 19 EPRPIAVGYV 29 D DTGFVRPD 38 SDAASQRHEP 48 RAPHIEQUOR 1 CSUSHRYF 9 YISVSRPGRG 19 SG EFHFDP 27 DGDEIFHVDH 37 AXX 40 EIVHRLEEPG 1 TKLEHVIIQA 11 EFYLH PDQ 19 SG EFHFDP 27 DGDEIFHVDH 37	
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DR1 6,	28 ERCIYNQEES 38 VIEDSDVGEY 48 KNY CENTER 5	
3000	172 LENGKETLQR	
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DR1 8,	87 ESFIVQRRVII	

Conserved residues Polymorphic residues

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ξì	6 VAL	CG1	47.07672	-36.30119	101.00126	Al	8	FF 10200005
€2	6 VAL	CG2		-37.46647	100.11057	A1	6	0.0000
63	6 VAL	C		-35.00190	103.02675		6	0.00000
64	6 VAL	0	-50.21018	-34.34041	102.87457		6	0.00000 0.00000
65	7 ILE	N	-48.11527	-34.52120	103.65431		7	0.00000
66	7 ILE	н	-47.35972	-35.13148	103.91409		'n	0.00000
67	7 ILE	CA	-48.09506	-33.08697	105.39701		ż	0.0000
68	7 ILE	CB	-48.69197	-32.86813 -33.71317	106.43001	Al	7	0.0000
69	7 ILE 7 ILE	CG2	-47.96322 -48.7/316	-31.39586	105.81727		7	0.00000
70	7 ILE 7 ILE	CG1	-49.74316	-31.20489	107.23523		7	0.0000
71 72	7 ILE	C	-46 693R1	-32.50114	103.87753	Al	7	0.00000
73	7 ILE	ō	-45.72315	-33.10109	104.32366		7	0.00000
74	8 ILE	N	-46.61414	-31.32789	103.23109		8	0.00000
75	8 ILE	н	-47.43598	-30.79777	103.00074		8	0.00000
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92	9 GIN	HE22	-42.66204	-25.53260	108.41320	Al	é	0.00000
93	9 GLN	С	42.00040	-27.79728 -28.56703	103.06902	λ1	9	0.0000
94	9 GLN	0 .	-47 01720	-26.84230	101.96835	Al	10	0.00000
95	10 ALA 10 ALA	N H	-42 B0016	-26.24022	101.81084		10	0.00000
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98 98	10 ALA	CB		-26.95993	99.69281	A1	10	0.00000
99	10 ALA	c	-40.41733	-25.14834	101.25674	Al	10	0.00000
100	10 ALA	o	-41.21138	-24.24680	101.50542	Al	10	0.00000 0.00000
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115	12 PHE	CG	-37.93682 -38.35291	-23.20899	96.17465		12	0.00000
116 217	12 PHE	CD1 CD2		-21.25353	95.22261	λl	12	0.00000
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123	13 TYR	<u>;</u> 2	-35.16677	-20.22034	99.15293 99.03348	E.Y	13	0.00000
124	13 TYR	H	-34.71888	-21.10856	33.02245	1		2.0000

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125	13 TYR	. CA	4.4465	0 -19.0289	1 99.57908	A1 13	0.00000
126	13 TYR	CB	3.7416	4 -19.3163	7 100.90471		0.00000
3.27	13 TYR	CG	-33.6322	9 -18.0484	4 101.71489		0.00000
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132	13 TYR			8 -15.7159			0.0000
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186	19 SER	н	-34.63228	-10.43617	101.06649	Al 19	0.00000
187	19 SEP	CA	-34.94474	-12.50616	100.83625	A1 19	0.00000 0.00000
188	19 SER	CB	-35.43672	-12.36114	102.28173	A1 19	5.00000

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189	19 SE	R OG	34.3706	64 -11 877	27 103.1084	1 11	19	0.00000
190					99 103.4025		19	0.00000
1 92			-35.9764	0 -13.297	76 100.0219	7 A1	19	0.00000
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193					37 100.5300		20	0.00000
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195	20 GL			3 -15.4753			20	0.00000
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207	21 GLU			6 -19.7836			21	0.00000
208 209	22 PHE 22 PHE		-42.3020	6 -19.3603 4 -18.7634			22 22	0.00000
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·211	22 PHE			-20.3120			22	0.00000
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213	22 PHE				6 105.4491		22	0.00000
214	22 PHE	CD2	-42.01384	-22.7696	6 104.14562	2 Al	22	0.00000
215	22 PHE	CEI	-40.30109	-22.3064	3 106.34948	Al	22	0.00000
216	22 PHE	CE2			0 105.04232		22	0.00000
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219	22 PHE	0			5 101.55111		22	0.00000
220	23 MET	И			101.32298		23 23	0.00000 0.00000
221 222	23 MET 23 MET	Н СА	-44,59132	-23.0327	9 101.30500 7 101.02967	ומי	23	0.00000
223	23 MET	CB		-22.73266			23	0.00000
224	23 MET	CG	-36.39952				23	0.00000
225	23 MET	SD	-46.67153	-23.10231		A1	23	0.00000
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228 229	23 MET	0			100.62115		23	000000
230	24 PHE 24 PHE	И Н			101.99622		24 24	0.00000 0.00000
231	24 PHE	CA			102.69060		24	0.00000
232	24 PHE	.CB			103.21139		24	0.00000
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238	24 PHE	Cz			106.72434		24	0.00000
239 240	24 PHE 24 PHE	Ç.,			101.85776		24 24	0.00000
241	25 ASP	О			101.74049		25	0.00000
242	25 ASP	н			102.20704		25	0.00000
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248	25 ASP	C	-49.18430				25	0.00000
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250 251	26 PHE	n H	-50.15819				26	0.00000
252	26 PHZ	CA	-50.56030				26	0.00000
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253	26 PHE	СВ	51.4237	8 -30.8309	3 102.82005) A1	26	0.00000
254	26 PHE	CG	-52.6035	6 -29.9203	8 102.57724		26	0.00000
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258	26 PHE				5 101.74257	' Al	26	0.00000
259	26 PHE	CZ		9 -28.1854			26	0.00000
260	26 PHE	С			100.51070		26	0.00000
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266	27 ASP	CG		5 -35.66101 0 -36.23164			27	0.00000
267 268	27 ASP 27 ASP	OD1 OD2	-52.95040) -36.23164) -36.23164	101.40272		27	0.00000
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273	28 GLY	CA		-33.86633			28	0.00000
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275	. 28 GLY	ō		.,~32.03342			28	0.00000
276	29 ASP	N		-31.96145			29	0.00000
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283	29 ASP	C		-29.64631	96.07233		29	0.00000
284	29 ASP	0		-28.59407			29 30	0.00000 0.00000
285	30 GLU	Ŋ		-29.81701	97.32405 97.90251		30	0.00000
286	30 GLU	H		-30.57647	97.86160	A1	30	0.00000
287	30 GLU	ÇA		-28.70871 -29.18152	98.54724	A1	30	0.00000
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290	30 GF0	CD		-29.50058	98.43828		30	0.00000
291	30 GLU	OE1		-30.47326	98.78969		30	0.00000
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297	31 ILE	Cλ		-25.52556		አ1	31	0.00000
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302	31 ILE	С			100.56211		31	0.00000
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304 305	32 PHE 32 PHE	N .			101.18020		32	0.00000
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300	32 PHE	CB	-50.76201	-25.18703	103.73391	Al	32	0.00000
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312	32 PHE	CE2	-52.82688	-27.19600	106.29575	A.1	32	0.00000
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315	32 PHE	0	-52.76581	-22.4:086	103.62205	λl	32	0.00000
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350
      39 LYS
               ΝZ
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./DR1_	_ಸಿದಾನ2 . ದಾಜ	>	wa Lop	25 14:58:	48 1993	7	
381	39 LYS	нг н	55.9564	5 -18.5458	1 100.3£313	λ1	3 0.00000
382	39 LYS		-55.3986	3 -17.2672	5 99.46103	λl	39 0.00000
3 نه 3	39 LYS	н н д з	-54.7248	6 -17.5192	1 100.90690	A1	39 0.00000
384	39 LYS				3 102.84(33		39 0.00000
385	39 LYS		-53.4335	1 -17.6751			39 0.00000
386	40 GLU		-53.4133	2 -19.7742	0 102.71885 1 103.42697	VI .	40 0.00000 40 0.00000
387 388	40 GLU		-53.1615	g -20.4369. 2 -20 14649	5 101.48663	A1	40 0.00000
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394	40 GLU		-55.46625	5 -20.74560	101.61501	Al (40 0.00000 40 0.00000
395	40 GLU		-56.42031	_20.18304	101.08479	V1 .	40 0.00000 41 0.00000
396 397	41 THR 41 THR		-50.56321	-21.01370	102.86321	Al	0.00000
398	41 THR		-56.84476	-22.55464	102.30963	Al 4	0.00000
399	41 THR	CB	-56.55011	-24.05830	102.16109	A1	0.00000
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401	41 THR	HG1		-25.59614			0.00000
402	41 THR	CG2		-24.67173			0.00000
.403	41 THR	C		-22.26510			61 0.00000 61 0.00000
404 405	41 THR 42 VAL	0		-21.93546 -22.41525			0.00000
406	42 VAL	и н		-22.81422			0.00000
407	42 VAL	CA		-22.03291			0.00000
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415	43 TRP	CA	-60.03340	-24.82113	106.34002		0.00000
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419	43 TRP	CE2	-60.47947	-27.49629	102.60102	Al 4	3 0.00000 3 0.00000
420 421	43 TRP 43 TRP	CE3 CD1	-58.30029 -62.11826	-21.24214 -26 B6040	104.12574 103.93459	A1 4	3 0.00000
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423	43 TRP	HE1	-62.50030	-27.49046	101.97106	A1 4	3 0.00000
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426	43 TRP	CH2	-58.24204	-28.07937	101.82597 107.16564		3 0.00000 3 0.00000
427 428	43 TRP 43 TRP	0	-59.07150	-25.043/3	108.02836	A1 4	3 0.00000
429	44 ARG	N			106.89519		4 0.00000
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435 436	44 ARG	NE	-52.18625	-23.26000	108.32672	A1 4	
437	44 ARG 44 ARG	HE CZ	-50.93149	-23.33650	107.86475	A1 4	
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439	44 ARG	нніі	-48.98076	-22.84912	108.24673	Al 4	
440	SAR FF		-50.05182	-22.52238	109.51721	hl 4	
441	44 ARG	NH2	-50.67215	-23.69024	106.69352	A1 4	
4 42 4 43	44 ARG	HH21	-49.73591	-23.96008	106.35139	A1 4	
444	44 ARG	∺∺22 C	-51.41769 -57.42751	-27,23910 -22,65857	108.38029	A) 4	
	11 770	~ .	24.72/21	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,		Ť	

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445	44 ARG	. 0	6.5306	4 -21.3920	1 108.88943	3 &1	44	0.0000
446	45 LEU	_		0 -21.5568			45	0.00000
447	45 LEU			6 -22.0638			45	0.00000
448	45 LEU			9 -20.2189			45	0.00000
449	45 LEU			7 -19.4651			45	0.00000
450	45 LEU		-59:1505	6 -18.9699	0 106.49579	Al	45	0.00000
451	45 LEU	CD1	-60:1453	4 -18.4190	6 105.47407	1 1 1	45	0.00000
452	45 LEO	CD2	-58:1043	1 -17.9228	9 106.87944	Y1	45	0.00000
453	45. Leu	С		8 -20.2010			45	0.00000
454	45 LEU	0	-59.5336	2 -19.1992	6 110.79199) Al	45	0.00000
455	46 GLT	N	-60.0891	7 -21.3602	4 110.48954	λì	46	0.00000
456	46 GLU	-	-60.1404	7 -22.1470	8 109.87631	λ1	4 6	0.00000
457	46 GLU		-60.5837	9 -21.4731	7 111.86481	Al	46	0.00000
458	46 GLU			7 -22.7151			46 46	0.00000 0.00000
459	46 GLU		-60.8088	1 -24.02385	111.51//2	N T	46	0.00000
460	46 GLU	CD OE1		4 -25.11830 7 -25.47222			46	0.00000
461 462	46 GLU 46 GLU	OE2		4 -25.6061			46	0.00000
463	46 GLU	C		3 -21.49323			46	0.00000
464	46 GLU	.0		-21.22226			46	0.00000
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.467	47 GLU	CA		-21.73864			47	0.00000
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469	47 GLU	ÇG		-23.50661			47	0.00000
470	47 GLU	CD	-55.55418	-24.58648	112.78701	λl	47	0.00000
471	47 GLU	QΞ1	-55.48886	-25.74409	112.37878	Al	47	∙0.00000
472	47 GLU	OE2	-55.29749	-24.27876	113.95212	λl	47	0.00000
473	47 GLU	С		-20.38151			47	0.00000
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475	48 PHE	N	-56.83807	-20.40716	115.24713	YJ	48	0.00000
476	48 PHE	R .	-56.92233	-21.28123	115.72305	λl	48	0.00000
477	48 PHE	CX	-56.75617	-19.13647	115.96362	Al	48	0.00000
478	48 PHE	CB	-57.20231	-19.34982	117.41420	YT.	48	0.00000 0.00000
479	48 PHE	CG	-58.41697	-18.49841	117.70829	VI	48 48	0.00000
480	48 PHE	CD1	-29.11175	-19.00320 -17.19966	117.44215	7.1	48	0.00000
401 482	48 PHE	CD2 CE1	-50.25455	-18.20670	117 71246	Al	48	0.00000
483	48 PHE	CE2		-16.40286			48	0.00000
484	48 PHE	CZ	-60 68202	-16.90878	118.24779	λl	48	0.00000
485	48 PHE	C	-55 39240	-18.46628	115.93777	λ1	48	0.00000
486	48 PHE	o	-54:35117	-19.07618	115.72587	Al	48	0.00000
487	49 GLY	N		-17.15361	116.19781		49	0.00000
488	49 GLY	В		-16.72101			49	0.00000
489	49 GLY	CV	-54.23420	-16.30520	116.13162	Al	49	0.00000
490	49 GLY	С	-52.92695	-16.86464	116.68056	A1	49	0.00000
491	49 GLY	0	-51.85815	-16.69886	116.10714	Y1	49	0.00000
492	50 ARG	N	-53.03915	-17.56211	117.81779	Y1	50	0.00000
493	SO ARG	H	-53.93660	-17.67284	118.24069	Al	50	0.00000
494	50 ARG	СУ	-51.84656	-16.16800	118.42224	Y1	50	0.00000
495	50 ARG	CB.	-52.27437	-18.92760	119.68340	WT.	50 50	0.00000
496	50 ARG	CG	-51.10260	-19.32736	120.38234	V.T	50	0.00000
497	50 ARG	CD	-54.55002	-20.14774 -20.44943	121.80047		50	0.00000
498 499	50 ARG	ΝE		-19.71957			50	0.00000
500	50 ARG	HE CZ		-21.63184	123.26571		50	0.00000
501	50 ARG	NH1		-21.83164			50	0.00000
502	50 ARG	8811	-49.05579	-22.04700	124,52612	A.1	50	0.00000
503	50 ARG	HH12	-48.49360	-21.13233	124.16453	۸ì	50	0.00000
504	50 ARG	NH2	-51.15926	-22.58629	123.10823	A.1	50	0.00000
505	50 ARG	HH21	-51.08073	-23.47414	123.56513	ሥጋ	50	0.00000
506	50 ARG	HH22	-51.95092	-22.42795	122.51852	Al	50	0.00000
507	50 ARG	С	-51.06703	-19.09561	117.49074	Al.	50	
508	50 ARG	0	-49.84240	-19.09133	117.41926	<i>k</i> .:	50	0.00000

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509	51 PHE	и :	8491	5 -19.8777	8 116.73839	A1	51	0.00000
510	51 PHE	-			7 116.76797		51	0.00000
513	SI PHE				6 115.73402		51	0.00000
512	51 PHE				3 115.15449		51	0.00000
513	51 PHE				4 114.02359		51	0.00000
514	51 PHE				114.29503		51	0.00000
515	51 PHE		-52.4578	8 -22.06484	112.69289	וג	51	0.0000
516	51 PHE		-51.72483	3 -24.74421	113.22969	A1	51	0.00000
517	51 PHE	CE2	-52.30778	3 -22.98082	2 111.63059	Al	51	0.00000
518	51 PHE	CZ	-51.94314	4 -24.31860	111.90000	λl	51	0.00000
519	51 PHE	С	-50.50268	-20.00696	114.69371	λl	51	0.00000
520	51 PHE	0	-49.31281	-20.20801	114.47961	Al	51	0.00000
521	52 ALA	.Ж	-51.20469	-19.03212	114.10202	Al	52	0.00000
522	52 ALA	H	-52.16823	3 -18.89764	114.33870	λ1	52	0.00000
523	52 ALA	CA	-50.54896	5 - 18.16486	113.11957	λl	52	0.00000
524	52 ALA	CB			112.62511		52	0.00000
525	52 ALA	С	-49.28257	-17.48933	113.62687	A1	52	0.00000
526	52 ALA	0			112.94510		52	0.00000
527	53 SER	N			114.88764		53	0.00000
528	53 SER		50.22411				53	0.00000
529	53 SER	CA	-48.18100	-16.47055	115.53815	Al	53 53	0.00000 0.00000
530	53 SER	CB	-48.58146	-16.02158	116.95383	W.I	53 53	0.00000
531	53 SER	- OG			117.52845		53 53	0.00000
532	53 SER	HG		-14.90357	118.42105 115.57426		53	0.00000
533 534	53 SER 53 SER	C	-46.99433	-17.42737 -17.11790	115.12568	VI.	53	0.00000
531 535	53 SER 54 PHE	O N.	-43.03403	-10 64770	116.08200	λl	54	0.00000
536	54 PHE	H.		-10.87361		Al	54	0.00000
537	54 PHP	CA	-46 18727	-19 64350	116.09999	A1	54	0.00000
538	54 PHE	CB.	-46 6954A	-20 99079	116.63413		54	0.00000
539	54 PHE	CG		-20.99411		λl	54	0.00000
540	54 PHE	CD1		-21.50316	118.65539		54	0.00000
541	S4 PHE	CD2		-20.53119	119.00621		54	0.00000
542	54 PHE	CEl		-21.55376			54	0.00000
543	54 PHE	CE2	-46.08993	-20.57973	120.40382	λl	54	0.00000
544	54 PHE	CZ	-47.30008	-21.09192	120.92307	A1	54	0.00000
545	54 PHE	С	-45.57270	-19.90650	114.73758	Al	54	0.00000
546	54 PHE	0	-44.36030	-19.93109	114.55246		54	0.00000
5 4 7	55 GLU	N		-20.08976	113.76558	λl	55	0.00000
5 48	55 GLU	H.			113.95516		55	0.00000
549	55 GLU	CA		-20.41545		λl	55	0.00000
550	55 GLU	CB		-20.81915		Al	55	0.00000
551	55 GLU	CC	-47.92567	-21.96126	112.19274	Al	55 55	0.00000
552	55 GLU	.CD			111.27677		55	0.00000
553 554	55 GLU 55 GLU	OE1 OE2	-49.13333	-23.33377	110.95047	Al	55	0.00000
555	55 GLU	C.	-45 10354	-19 30188	111.76318	λl	55	0.00000
55.6	55 GLV	0:	-44 12678	-19 50387	111.19343	λ1	55	0.00000
557	56 ALA	N	-45 73650	-18 08673	111.89532	A1	56	0.00000
558	56 ALA	H	-46.62336	-17.96765	112.34890	Al	56	0.00000
559	56 ALA	CA	-45.00414	-16.92733	111.38640	Al	56	0.00000
560	56 ALA	CB	-45.80074	-15.63871	111.59969	Al	56	0.00000
561	56 ALA	c	-43.63772	-16.77849	112.02791	Al	56	0.00000
562	56 ALA	0	-42.62065	-16.60634	111.36878	Al	56	0.00000
563	57 GLN	N	-43.63088	-16.91454	113.35929	Al	. 57	0.00000
564	57 GLN	H	-44.48826	-17.03983	113.86832	λl	57	0.00000
565	57 GLN	CΛ	-42.35063	-16.88200		λl	57	0.00000
566	57 GLN	CB	-42.61967	-17.03049	115.56930	Al	57	0.00000
567	57 GLN	CG	-61.40659	-15.75355	116.45895	A.1	57	0.00000
566	57 GLN	CD	-41.80070	-16.94589	117.90800	A.1	57 57	0.00000 0.00000
569	57 GLN	OEl	-42.29879	-16.06206	118.58910	A1	57 57	0.00000
570	57 GLN	NE2	-41.55425	-10.16011	110.36366	E1	57 57	0.00000
571	57 GLN	HE21	-41.14709	-18.87175	110 33353	ום 1	57	0.00000
572	57 GLN	HEZZ	-41.78115	-15.36/28		***		0.0000

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573	57 GLN	•	1 3621	3 ~17 04055	113.59025	λl	57	0.00000
574	57 GLN		~40.1897	1 -17.68051	113.34773	Al	57	0.00000
373	58 GLY				113.41319		8 ئ	0.00000
576	58 GLY		-42.8477	1 -19.33514	113.64847	λl	58	0.0000
577	58 GLY		-41.05379	9 -20.23393	112.87374	A1	58	0.00000
578	58 GLY	С	-40.45534	4 -19.92781	111.50957	Al	58	0.00000
579	58 GLY	0	-39.2532	-20.01021	111.27849	λl	5 B	0.00000
580	59 ALA	N	-41.34654	-19.52227	110.60018	NI NI	59 59	0.00000
581	59 ALA	Н	-42.32072	2 -19.46/62	110.83485	V1	59	0.00000
582	59 ALA	CA	-40.88666	-19.14337	109.20101	A I	59	0.00000
583	59 ALA 59 ALA	CB C	-32.0/124	1 -18.74000 1 -18 02257	109.26073	Al	59	0.00000
584 585	59 ALA	0	-39.86090 -38.85187	-18.0410; -18.04905	108.56603	A1	59	0.00000
585 586	60 LEU	N.	-40.12979	-17.02925	110.11113	Al	60	0.00000
587	60 LEU	н	-40.96678	-17.03715	110.66448	Al	60	0.00000
588	60 LEU	CA	-39.17026	-15.93079	110.21454	λl	60	0.00000
589	50 LEU	CB	-39,82974	-14.72108	110.88234	Al	60	0.00000
590	60 LEU	CG	-41.00342	-14.16448	110.06267	Al	60	0.00000
591	60 LEU	CD1	-41.72077	-13.05637	110.83175	A1	60	0.00000
592	60 LEU		40.54968	-13.69403	108.67852	VI	60 60	0.00000
593	60 LEU	C	-37.86300	16.28687	110.90549	7.7	60	0.00000
594	60 LEU	0	-30.81366	12 20628	110.64266 .111.76650	A 1	61	0.00000
595 596	61 ALA 61 ALA	И Н	-30.92346	-17.30828	112.04737	Al	61	0.00000
597	61 ALA	n CA	-36.66060	-17.70200	112.25036	A1	62	0.00000
59E	61 ALA	CB	-36 90091	-18.87147	113.37402	Al	61	0.00000
599	61 ALA	c	-35.86652	-18.52303	111.13575	A1	61	0.00000
600	61 ALA	. 0	-34.67753		110.93578	Al	61	0.00000
601	62 ASN	Ň	-36.59182	-19.33811	110.35468	Al	62	0.00000
602	62 ASN	H	-37.55651	-19.52444	110.56458	A2	62	0.00000
603	62 ASN	CA	-35.93048	-19.97053	109.20954	Al	62	0.00000
604	62 ASN	ĊВ	-36.90608	-20.83397	108.41185	Al	62	0.00000
605	62 ASN	CG	-36.14296	-21.97501	107.76767	A1	62	0.00000 0.00000
606	62 ASN	OD1	-35.90083	-23.00051	108.38915	A.L	62 62	0.00000
607	62 ASN	ND2	-35.81296	-21.80385	106.49294	Y I	62	0.00000
608	62 ASN	HD21	-35.71613	-20.90910	105.92778	2.1	62	0.00000
609 610	62 ASN 62 ASN	НФ22 С		-18.97317			62	0.00000
611	62 ASN	0		-19.05772		Al	62	0.00000
612	63 ILE	N	-36.07385	-17.96130	107.91224		63	0.00000
613	63 ILE	H	-37.03805	-17.97906	108.18770	Y]	63	0.00000
614	63 ILE	CA	-35.60960	-16.86395	107.05550	λl	63	0.00000
615	63 ILE	CB	-36.79680	-15.88630	106.84927	Al	63	0.00000 0.00000
616	63 ILE	CG2	-36.71800	-14.58751	107.66300	Al Il	63 63	0.00000
617	63 ILE	CG1	-37.00443	-15.60068	105.06416		63	0.00000
618 619	63 ILE 63 ILE	Ċ C	-36.27101	-16 14412	107.48562	λl	63	0.0000
620	63 ILE	0	-33.67028	-15.43835	106.72047	አነ	63	0.00000
621	64 ALA	N	-33.97867	-16.34078	108.76481	Al	64	0.0000
622	64 ALA	н	-34.55914	-16.88095	109.37800	አነ	64	0.00000
623	64 ALA	CA	-32.68252	-15.86370	109.23001	Al	64	0.00000
624	64 ALA	CB	-32.78414	-15.37252	110.67448	Al	64	0.00000
625	64 ALA	С	-31.59324	-16.91956	109.13883	Al	64	0.00000
626	64 ALA	0	-30.45701	-16.65389	108.75412	Al	64	0.00000
627	65 VXI	N	-31.96250	-18.15367	109.50633	F)	65 65	0.00000
628	65 VAL	H	-32.92093		109.71401		65	0.00000
629	65 VAL	CA	-30.91507 -31.28412	-19.17596	110 77710	21	65	0.00000
630 53)	65 VAL	CB CG1	-31.28412	-19 97575	111.77524	Al	65	0.0000
631 632	65 VAL 65 VAL	CG1 CG2	-32.45603	-21.23819	109.79675	A2	65	0.00000
633	65 VAL	C C	-30,45713	-19.56758	108.08237	Y.	65	0.0000
634	65 VAL	ō	-29.26568	-19.71596	107.82437	W.	65	0.00000
635	66 ASP	N	-31.42136	-19.66037	107.15605	Al	65	0.00000
636	66 ASP	H	-32.38553	-19.49723	107.38932	ř. 1	6 €	0.00000

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63.7	66 ASP	CA	1.0485	8 -19.93210	6 105.76795 A	66	0.0000
638	66 ASP				6 104.93007 A	66	0.00000
533	S6 ASP				105.08208 A		0.00000
640	66 ASP				105.20729 3		0.00000
641	66 ASP			6 -19.6546			0.00000
642	66 ASP		-30.18170	8 -18.8344;	9 105,15564 ك 2 104,50705 كى		0.00000
643 644	66 ASP 67 LYS		-29.16484 -30.56424	2 -19.00072 1 -17 5991	5 105.48017 A	67	0.00000
645	67 LYS		-31 49536	-17.48655	105.84055 A	67	0.00000
646	67 LYS	CV.		-16.43204			0.00000
647	67 LYS	СВ	-30.39071	-15.23196	3 105.87104 A		0.00000
648	67 LYS	CG	-29.79159	-13.86177	1 105.56853 AI	. 67	0.00000
649	67 LYS	CD	-30.51506	-12.79056	106.38164 31	. 67	0.00000
650	67 LYS	CE	-29.96446	-11.38645	106.14720 Al	67	0.00000
651	67 LYS	NZ		-10.44891	. 107.01869 A1	67 67	0.00000
652	67 LYS	HZ1 HZ2	-30.33262		106.83909 21	67	0.00000
653 654	67 LYS 67 LYS	H23	-31.70143	-10.71804	108.01156 كل	67	0.00000
655	67 LYS	c	~20.28117	-16.58093	105.64383 A1	67	0.00000
656	67 LYS	. 0	27.33559	16.44078	104.87661 Al	67	0.00000
657	68 ALA	N	-28.12520	-16.92373	106.92795 Al	68	0.00000
658	68 YTY	H	-28.92119	-17.00935	107.53692 31	68	0.00000
659	68 ALA	CA	-2676352	17.18143	107.40958 Al	68 68	0.00000
660 661	68 ALA	CB	-26.77377	-17.44846	108.91534 Al 106.70618 Al	68	0.00000
662	68 ALA 68 ALA	С 0		-18.33297		68	0.00000
663	69 ASN	N	-26.87877			69	0.00000
.664	69 ASH	н	-27.84416	-19.34421	106.72158 Al	69	0.00000
665	69 ASN	CA	-26.32826	-20.54731	105.77098 Al	69	0.00000
666	69 ASN	CB	-27.33794	-21.70567	105.74618 31	69	0.00000
667	69 ASN	CG	-27.75534	-22.20215	107.12937 או	69	0.00000
668	69 ASN	OD1	-28.81753	-22.77967	107.30600 A1	69 69	0.00000 0.00000
669	69 ASN	ND2	-26.90880 -26.02949		108.13718 A1 108.03476 A1	69	0.00000
670 671	69 ASN 69 ASN	HD21	-20.02949	-22 29754	109.04652 Al	69	0.00000
672	69 ASN	C	-25.83413	-20.26827	104.36379 A1	69	0.00000
673	69 ASN	ō	-24.88019	-20.87816	103.89106 Al	69	0.0000
674	70 LEU	И	-26.46696	-19.27268	103.71664 Al	70	0.00000
675	70 LEU	H	-27.27121	-18.83110	104.12686 A1	70 70	0.00000 0.00000
676	70 LEU	CA	-25.93555	-18.80513	102.42930 Al	70	0.00000
677 678	70 LEU 70 LEU	CB	-26.70466	-17 87907	101.93156 A1 101.32608 A1	70	0.00000
679	70 LEU 70 LEU	CG CD1	-28 90878	-16.60756	101.23109 A1	70	0.0000
680	70 LEU	· CD2	-27.93286	-18.56192	99.96378 A1	70	0.00000
681	70 LEU	С	-24.47328	-18.42736	102.51389 Al	70	0.00000
682	70 LEU	0	-23.64160	-18.86456	101.72791 AL	70	0.00000
683	71 GLU	N	-24.17065	-17.62592	103.54240 A1	71 71	0.00000 0.00000
684	71 GLU	H	-24.87529	-17.33320	104.19342 Al	71	0.00000
685 686	71 GLU 71 GLU	CA CB	-22.77384	-16 23884	103.73624 A1 104.88750 A1	71	0.00000
687	71 GLU	CG	-21.33647	-15.50613	104.93364 Al	71	0.00000
688	71 GLU	CD	-21.30052	-14.53829	106.10023 A1	71	0.00000
689	71 GLU	OE1	-20.27115	-14.47665	106.76970 Al	71	0.00000
690	71 GLU	OE2	-22.29376	-13.85012	106.33423 Al	71	0.00000
691	71 GLU	С	-21.86369	-18.43808	103.97868 A1	71 71	0.00000 0.00000
692	71 GLU	0 ,	-20.81243	-18.61050	103.36820 Al 104.86364 Al	72	0.00000
693	72 ILE 72 ILE	N U	-22.34609	-19 12122	105.33200 Å	72	0.00000
694 695	72 ILE 72 ILE	H AO	-23.21301	-20.53142	105.13649 Al	72	0.00000
696	72 ILE	CE	-22.29516	-21.40405	106.17923 21	72	0.00000
697	72 ILE	CG2	-21.51796	-22.68646	106.50604 ሕነ	72	0.00000
696	72 ILE	CG1	-22.55172	-20.59465	107.45409 A1	72	0.00000
699	72 ILE	CD	-23.34520	-21.36378	108.51180 Al	72 72	0.00000 0.00000
730	72 ILE	С	-21.22106	-21.35113	103.89490 A1	• 2	5.0000

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701	72 ILE	. 0	20.06679 -21.68162 103.64178 A1	72	0.00000
702	_		-22.24406 -21.67127 103.09738 Al	73	0.00000
763		н	-23.17806 -21.34412 103.27303 Al	73	0.00000
704	73 MET	CA	-21.88577 -22.51146 101.95532 A1	73	0.00000
705			-23.02917 -23.44601 101.56487 A1	73	0.00000
706			-23.30745 -24.44882 102.69030 A1	73	0.00000
707	73 MET		-24.39480 -25.79652 102.19686 A1	73	0.00000
708	73 MET		-24.58745 -26.56731 103.81212 Al	73	0.00000
709	73 MET		-21.30840 -21.79430 100.75010 A1	73	0.00000
710	73 MET	_	-20.54740 -22.36938 99.97894 A1	73	0.00000
711	74 THR		-21.60342 -20.49214 100.63510 A1	74	0.00000
712	74 THR 74 THR		-22.28955 -20.04403 101.21510 A1	74	0.00000
713 714	74 THR 74 THR		-20.82593 -19.73833 99.64704 A1 -21.46299 -18.35496 99.36040 A1	74 74	0.00000 0.00000
715	74 THR		-21.46299 -10.33496 99.36040 AT -21.01301 -17.84761 98.09669 A1	74	0.00000
716	74 THR	HG1	-20.04699 -17.88381 98.02785 A1	74	0.00000
717	74 TIIR	CG2	-21.20740 -17.31089 100.44667 A1	74	3.00000
718	. 74 THR	C	-19.35170 -19.61359 100.02467 A1	74	0.00000
719	74 THR	Ó	-18.48554 -19.49348 99.16750 A1	74	0.00000
- 720	75 LYS	·N	· -19.08538 -19.69260-101.33717- A1	75	0.00000
721	75 LYS	н	-19.81075 -19.65220 102.02321 A1	75	0.00000
722	75 LYS	CX	-17.69833 -19.83955 101.77078 Al	75	0.00000
723	7,5 LYS	CB	-17.61408 -19.57287 103.27797 A1	75	0.00000
724	75 LYS	CG	-16.20208 -19.66153 103.86107 A1	75	0.00000
725	75 LYS	CD	-16.21091 -19.57437 105.38567 A1	75	0.00000
726	75 LYS	CE	-14.81504 -19.74395 105.98417 A1	75	0.00000
727	75 LYS	NZ	-14.91179 -29.69891 107.44939 A1	75	0.00000
728 729	75 LYS	HZ1	-13.96820 -19.83508 107.86470 A1	75 75	0.00000
730	75 LYS 75 LYS	H22 H23	-15.29159 -18.77551 107.74094 A1 -15.54844 -20.45311 107.77639 A1	75 75	0.00000
731	75 LYS	C.	-17.14118 -21.21778 101.45102 A1	75	0.00000
732	75 LYS	o.	-16.11623 -21.36022 100.79660 A1	75	000000
733	76 ARG	N	-17.86151 -22.25466 101.90990 A1	76	0.00000
734	76 ARG	H	-18.69512 -22.09296 102.44436 A1	76	0.00000
735	75 ARG	CA	-17.35520 -23.61422 101.67525 A1	76	0.00000
736	76 ARG	CB	-18.33654 -24.68494 102.17145 Al	76	0.00000
737	76 ARG	CG	-18.82789 -24.66219 103.62361 A1	76	0.00000
738	76 ARG	CD	-19.55383 -25.98411 103.91795 A1	76	0.00000
739	76 ARG	NE,	-20.36155 -25.99642 105.14334 A1	76 76	0.00000 0.00000
740 741	76 ARG 76 ARG	E S	-21.34527 -25.85364 105.02467 A1	76	0.00000
742	76 ARG	CZ NH1	-19.85069 -26.28686 106.34915 A1 -20.67426 -26.45770 107.38550 A1	76	0.00000
743	76 ARG	HHII	-20.32838 -26.64144 108.30551 Al	76	0.00000
744	76 ARG	HH12	-21.67206 -26.42100 107.26619 A1	76	0.00000
745	76 ARG		-18.53304 -26.41209 106.51158 A1	76	0.00000
746	76 ARG		-18.12377 -26.62371 107.39854 A1	76	0.00000
747	76 ARG		-17.92942 -26.29542 105.72237 A1	76	0.00000
748	76 ARG	С	-17.06072 -23.91761 100.20901 A1	76	0.0000
749	76 ARG	0	-16.02715 -24.45819 99.83552 A1	76	0.00000
750	77 SER	N 	-18.01091 -23.50462 99.36716 A1	77	0.00000
751 752	77. SER	H.	-18.84286 -23.05363 99.69917 Al	77 77	0.00000 0.00000
753	•	CA	-17.80506 -23.71642 ,97.93702 A1 -19.13837 -24.15898 97.32351 A1	77	0.00000
754		CB OG	-19.13837 -24.15898 97.32351 A1 -18.92403 -25.08061 96.24728 A1	77	0.00000
7 5 5	77 SER	HG	-18.99574 -24.62840 95.39572 Al	77	0.00000
756		C	-17.22854 -22.51234 97.20137 A1	77	0.00000
757			-17.39865 -22.33993 96.00106 A1	77	0.00000
758	_		-16.52714 -21.66944 97.97646 A1	78	0.00000
759		H	-16.44843 -21.84082 98.96113 Al	78	0.00000
760			-15.74721 -20.53857 97.45264 A1	78	0.00000
761			-14.33390 -21.01377 97.09195 A1	7 E	0.00000
762			-13.52905 -21.21984 98.36114 A1	70 76	0.00000
763			-12.89966 -20.31632 98.89264 A1	7 E 7 S	0.00000 0.00000
754	78 ASK 1	ND2	-13,55264 -22,45457 98.85326 A1	, 5	0.0000

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765	78 ASN	нр2:	1 4.08412	2 -23.19170	98.43678 Al	78	0.00000
766	78 ASN	HD2		1 -22.65362	99.68493 Al		0.00000
767	78 ASN			9 -19.72172	96.31352 A1	78	0.00000
768	78 ASN			3 -19.42979	95.30178 Al	78	0.00000
769	79 TYR			-19.32597	96.52354 A1	79	0.00000
770	79 TYR			-19.55771	97.38252 Al	79 79	0.00000
771	79 TYR 79 TYR			3 -18.48918 3 -18.92649	95.49858 A1 95.21260 A1	79	0.00000
772 773	79 TYR			-20.32764	94.65360 Al	79	0.00000
774	79 TYR		-20 64246	-21.23278	95.25950 A1	79	0.00000
775	79 TYR		-20.76017	-22.54189	94.75217 Al	79	0.00000
776	79 TYR			-20.72112	93.54108 A1	79	0.00000
777	79 TYR			-22.03339	93.03642 A1	79	0.00000
778	79 TYR			-22.93822	93.64752 Al	79	0.00000
779	79 TYR	ОН		-24.23438	93.18578 Al	79	0.00000
780	79 TYR	нн	-19.67682	-24.30670	92.31232 A1	79	0.00000
781	79 TYR	C		-17.01261	95.83719 A1	79	0.00000
782	79 TYR	0		-16.51273	96.76683 Al	79	0.00000
783	80 THR	N		-16.33235	95.00452 A1	80 80	0.00000
784 785	80 THR	H		-16.79509	94.24288 A1 95.20513 A1	80	0.00000
786	80 THR 80 THR	CA CB		-14.92660 -14.87711	96.18098 Al	80	0.00000
787	80 THR	OG1		.,-13.52440	96.43537 Al	80	0.00000
788	80 THR	HGl		-13.50159	96.80560 A1	80	0.00000
789	80 THR	CG2.			95.72104 A1	80	0.00000
790	80 THR	С	-16.77512		93.83840 Al	80	0.00000
791	80 THR	O	-16.16879	-14.97644	92.99811 Al	80	0.00000
792	81 PRO	N	-17.23243	-13.08096	93.61451 A1	81	0.00000
793	81 FRO	CD	-18.04687		94.50529 A1	81	0.00000
794	81 PRO	CA	-16.95963		92.32255 A1	81	0.00000
795	81 PRO	СВ	-18.08102		92.28996 A1	81	0.00000 0.00000
796	81 PRO	ÇG	-18.22970		93.74456 X1 92.25040 A1	81 81	0.00000
797 798	81 PRO 81 PRO	C	-15.57247 -15.41926		92.16776 A1	81	Ú.00000
799	82 ILE	Р. И.	-14.55883		92.27154 A1	82	0.00000
800	82 ILE	H.	-14.69364	-13.66392	92.31920 A1	82	0.00000
801	82 ILE	CA		-12.18130	92.13921 A1	82	0.00000
802	82 ILE	СВ		-11.87598	93.53959 A1	82	0.0000
803	82 ILE	CG2	-12.41140	-13.13773	94.38674 Al	82	0.00000
804	82 ILE	CG1	T T,	-11.05228	93.44331 Al	82	0.00000
805	82 ILE	CD	-10.76554		94.80896 A1	82 82	0.00000 0.00000
8 <u>0</u> 6 807	82 ILE 82 ILE	C		-13.19907 -12.81747	91.37376 A1 90.75062 A1	82	0.00000
808	82 ILE	OCT1	-12.72556		91.38671 A1	82	0.00000
809	83 CLY	N	-17.53322	-0.31236	94.99084 B1	1	0.0000
810	83 GLY	HT1	-17.21994	0.44323	94.35235 B1	1	0.00000
811	83 GLY	HT2	-16.86357	-1.12219	94.96444 Bl	1	0.00000
812	83 GLY	нтЗ	-17.61098	0.01920	95.97150 B1	1	0.00000
813	83 GLY	CA		-0.91116	94.55151 B1	1	0.00000
814	83 GLY		-18.52573	-2.38203	94.66351 Bl	1	0.00000
815	83 GLY			2.70920	94.84086 B1	1 2	0.00000 0.00000
816 817	84 ASP		-19.57260	-3.20239	94.59303 B1 94.35600 B1	2	0.00000
818	84 ASP 84 ASP		-20.49658 -19.43900	-2.89510 -4.63200	94.86181 B1	2	0.00000
819	84 ASP		-19.44643	-4.83356	96.38475 B1	2	0.00000
82D	84 ASP		-18.89301	-6.19619	96.73815 Bl	2	0.00000
821	84 ASP		-17.69451	-6.29764	96.98109 B1	2	0.0000
822	84 ASP	OD2	-19.66566	-7.14958	96.75715 B1	2	0.00000
923	84 ASP		-20.62786	-5.31072	94.19848 Bl	2	0.00000
824	84 ASP		-21.46903	-4.61697	93.63479 B1	2	0.00000
925	85 THR		-20.67796	-6.64606	94.24891 Bl	3 3	0.00000
926 223	85 THR		-20.04362	-7.15926	94.84032 B1 93.59154 Bi	3	0.00000
827	65 THR		-21.75257 -21.58903	-7.39367 -7.31950	93.39134 E1 92.05122 E1	3	0.00000
328	95 THR	C5	-22.26903	-1.51350	,.,.,,,,	-	

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829	85 TH	R OG1	32.7626	5 -7.82486	91.39964 B1	3	0.00000
830			-22.7343				0.00000
83:			-20.3296				0.00000
832	85 TH		-21.7429			3	0.00000
833	85 TH		-20.7645			3	0.00000
834	86 ARC		-22.8649			4	0.00000
835	86 ARC		-23.6240			4	0.00000
836	86 ARC			0 -10.87552		4	0.00000
837	86 ARC			8 -10.79083		4	0.00000
838	86 ARC			3 -11.65365		4	0.00000
839	86 ARG			6 -13.10294		4	0.00000
840	86 ARG			-14.03911		4	0.00000
841	86 ARG			3 -14.68836		4	0.00000
842	86 ARG	CZ		-14.05016		4	0.00000
843	86 ARG	NHl		-14.85617		4	0.00000
. 844	86 ARG		-19.73815	-14.87872	99.26299 B1	4	0.00000
845	86 ARG	HH12	-20.05194	-15.47498	97.69021 B1	4	0.00000
846	36 ARG			-13.26636	99.42563 B1	4	0.00000
847	86 ARG				100.26897 B1	4	0.00000
848		HH22			- 99.36831 Bl	4	0.00000
849	86 ARG			÷11.82865	93.57905 B1	4	0.00000
850	86 ARG	0		-11.52883		4	0.00000
851	87 PRO	N		-12.98532	93.27325 B1	5	0.00000
852	87 PRO	CD		-13.34562	93.54376 B1	5	0.00000
B53	87 PRO	CA		-14.02757	92.53087 B1	5	0.00000
854 855	87 PRO	CB		-15.13066	92.39999 B1	5	0.00000
856	87 PRO	CG		-14.43376	92.52460 B1	5	0.00000
857	87 PRO 87 PRO	C		-14.54496	93.21975 B1	5 5	0.00000
858	87 PRO 88 ARG	0		-14.39544	94.41838 B1	6	0.00000
859	88 ARG	N H		-15.35588	92.39123 B1 91.44433 B1	6	0.00000
860	98 ARG	CA		-15.62675	92.90235 B1	6	0.00000
861	BP ARG	CB		-15.39805	91.81653 B1	6	0.00000
862	88 ARG			-15.55859	92.35844 B1	6	0.00000
B 63	88 ARG		-30.79150		91.49236 B1	6	0.00000
864	88 ARG		-31.92741		92.34607 B1	6	0.00000
8 6 5	88 ARG		-31.76512		93.33788 B1	6	0.00000
866	88 ARG		-33.08381		91.86193 B1	6	0.00000
8 67	88 ARG		-34.06565		92.71322 B1	6	0.00000
8 6 B	88 ARG	нніі	-34.95441	-13.48197	92.40144 B1	6	0.00000
869	88 ARG	HH12	-33.91853	-13.92533	93.69838 B1	6	0.00000
870	88 ARG	NH2	-33.26623	-13.97383	90.54974 B1	6	0.00000
871	88 ARG		-34,12737		90.16725 Bl	6	0.0000
872	88 ARG		-32.52080		89.92214 B1	6	0.00000
873	BB ARG		-27.23157		93.36366 B1	6	0.00000
874 875	BB ARG		-26.89591		92.63574 B1	6	0.00000
876	89 PHE		-27.62757		94.62546 B1	7	0.00000
877	89 PHE		-27.92346		95.18173 B1	7 7	0.00000
878	89 PHE		-27.64368 -26.56379		95.17433 B1 96.25427 B1	7	0.00000
879	89 PHE		-25.20774			7	0.00000
880	89 PHE		-24.74915	•	95.61829 B1 94.64858 B1	7	0.00000
881	89 PHE		-24.74915		95.96102 B1	7	0.00000
882	89 PHE		-23.51045		94.00961 B1	7	0.00000
883	89 PHS		23.18767		95.32384 B1	,	0.00000
884	89 PHE		22.73800		94.34976 B1	ż	0.00000
885	89 PHE		28.99632		95.72084 B1	7	0.00000
886	89 PHE			-18.12429	95.94786 B1	ז	0.00000
887			29.15750		95.87791 81	8	0.00000
885	90 LEU		26.38720 -		95.76376 B1	8	0.00000
889			30.48975 -		96.14113 Bi	8	0.00000
890			31.01265 -		94.84981 B1	8	0.00000
691				-21.16892	94.43614 B1	3	0.00000
992	90 LEU	CD1 -	32.81137 -	-21.96230	93.19113 21	3	0.00000

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8 9 3	90 LE	CD2	3,4749	8 -21.4069	9 95.55222	В1	8	0.00000
894	90 LEU			7 -21.9028			8	0.00000
895	90 LEU	0		6 -22.8920			8	0.00000
896	91 TRP			3 -21.7108			9.	0.00000
897	91 TRP			7 -20.8522			9 9	0.00000
898	91 TRP			7 -22.8604	3 99.06700 9 100.54418		9	0.00000
899 900	91 TRP 91 TRP				0 101.18157		9	0.00000
901	91 TRP			3 -22.7928			9	0.00000
902	91 TRP		-28.3007	4 -21.9654	6 101.98847		9	0.00000
903	91 TRP			8 -24.0032		B1	9	0.00000
904	91 TRP		-30.2650	0 -20.9642	7 101.96112	B1	9	0.00000
905	91 TRP	NEI	-28.9981	0 -20.8704	7 102.44030	B1	9	0.00000
906	91 TRP	HE1	-28.6856	6 -20.1561	5 103.04137	B1	9	0.00000
907	91 TRP	CZ2	-26.9776	7 -22.40520	102.16578	B1	9	0.00000
908	91: TRP	CZ3		-24.40402			9 9	0.00000
909	91 TRP	CH2		-23.61859			9	0.00000
910 911	91 TRP	C		5 -23.39534 5 -22.66026			9	0.00000
912	92 GLN		33.02847				10	0.00000
913	92 GLN	н	-32.20362	-25.28365	98.76579		10	0.00000
914	92 GLN	CA		-25.34751			10	0.00000
915	92 GLN	CB		,-26.04119		Bl	10	0.00000
916	92 GLN	CG	-34.43287	-25.05724	96.13126		10	0.00000
917	92 GLN	CD		-25.79390			10	0.00000
918	92 GLN	OEl		-25.88838			10	0.00000
919	92 CTN	NE2		-26.32196			10	0.00000
920	92 GLN	HE21		-26.21874			10	0.00000
921	92. GLN	HE22		-26.83019			10 10	0.00000
922	92 GLN	C		-26.35610	100.00960		10	0.00000
923 924	92 GLN 93 LEU	0. N.	-35.49484			Bl	11	0.00000
925	93 TEA	н.	-36.24432	-25.74963	100.29272		11	0.00000
926	93 LEU	CΛ	-35.59559	-27.43909	101.48633	91	11	0.00000
927	93 LEU	CB	-35.41178	-26.81504	102.88181	B1	11	0.00000
928	93 LEU	CG.	-34.83577	-27.70766	103.99848	Bl	11	0.00000
929	93 LEU	CD1	-34.33031	-26.83824	105.14843	Bl	11	0.00000
930	93 LEU	CD2	-35.82812	-28.74167	104.52805	2.1 2.1	11 11	0.00000 0.00000
931	93 LEU	Ċ	-36.93596	-28.11112	101.32695	B.	11	0.00000
932 933	93 LEU 94 LYS	И. О	-37.91092	-20 41719	101.56549	Bl	12	0.00000
934	94 LYS	н	-36.10205	-29.89866	101.79736	B1	12	0.00000
935	94 LYS	CA	-38.21063	-30.14121	101.49214	Bl	12	0.00000
936	94 LYS	CB	-38.26049	-30.88530	100.13839	Bl	12	0.00000
937	94 LYS	CG	-39.43761	-31.85442	100.07544	Bl	12	0.00000
938	94 LYS	CD		-32.44832	98.75113	RI	12	0.00000
939	94 LYS	CE		-33.40442	99.23924		12	0.00000
940	94 LYS	NZ		-33.90963	98.26846		12 12	0.00000 0.00000
941 942	94 LYS	H21		-34.41995 -33.14263	98.86351 97.78947		12	0.00000
943	94 LYS	HZ2 HZ3	-42.49249		97.59540		12	0.00000
944	94 LYS	C	-38.34910	-31.08699	102.67642		12	0.00000
945	94 LYS	ō	-37.54720	-31.99770	102.85849	Bl	12	0.00000
946	95 PHE	N	-39.40676	-30.85929	103.46609	Bl	13	0.00000
947	95 PHE	н	-39.99321	-30.05201	103.34519	Bl	13	0.00000
948	95 PILC	CA	-39.73794	-31.84982	104.49053	B1	13	0.00000
949	95 PHE	CB	-39.13251	-31.49228	105.86481	E1	13	0.00000
950	95 PHE	CC	-39.62104	-30.20819	106.49960	51 13	13 13	0.00000 0.00000
951	95 PHE	CD1	-39.04917	-28.96886	106.12996	BJ DT	13	0.00000
952 953	95 PHE	CD2	-40.00008	-20.40519	107.51173	B1	13	0.00000
953 954	95 PHE 95 PKE	CE1 CE2	-37.3330/ -41 01477	-29 08019	108.16313	B1	13	0.00000
955	95 PHE	CZ	-40.43557	-27.84435	107.79605	<u>۲</u>	13	0.00000
956	95 PHE	c	-41.22005	-32.16799	104.57423	Еì	13	0.00000

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95.7	95 PHE	•	08591	-31.31332	104.41691	Bl	1'3	-05-00000
958	96 GLU	0	47812	-33.46734	104.75822	Bl	14	0.00000
9::9	76 GLU	н	-40.74093	-34.09871	105.01948		14	0.00000
960	96 GLU	CA	-42.80002	-33.99059	104.40078	Bl	14	0.00000
961	96 GLU	CB	-42.75085	-34.25375	102.88321		14 14	0.00000 0.00000
962	96 GIU	CG	-43.92070	~34.87965	102.11791 100.64018		14	0.00000
963	96 GLU	CD.	-43.55505	-34.92496 -34.15253	99.84297	Bl	14	0.00000
964	96 CM	OEl	-44.08203	-35.68559	100.23154		14	0.00000
965 966	96 GLU	OE2 C	-43.13129		105.17925	B1	14	0.00000
967	96 GLÜ	0	-42.27510		105.37879	Bl	14	0.00000
968	97 CYS	N	-44 30621	-35.34431	105.62501	B1	15	0.00000
969	97 CYS	н	-45.05281	-34.60566	105.44206	B1	15	0.00000
970	97 CYS	CA	-44.78990	-36.60747	106.25700	Bl	15	0.00000 0.00000
971	97 CYS	CB	-45.09404		107.76554	BI	15 15	0.00000
972	97 CYS	SG	-46.49873		108.42288 105.55706	ומ דם	15	0.00000
973	97 CYS	C	-45.89248		103.33700	B1	15	0.00000
974	97 CYS	0	-46.75253	-38.69951	105.71784	Bl	16	0.00000
975	98 HIS	N	-45.80356 -45.06854		106.26371	B1	16	0.00000
976	98-HIS	·H	-46.75592	-39.59530		B 1	16	0.0000
977 978	98 HIS 98 HIS	CA CB	-46.00867	-40.61029	104.19437	Bl	16	0.00000
979	98 HIS	CG	-45.38253	-39.98357	102.97364	Bl	16	0.00000
980	98 HIS	ND1	-45.88967	-40.13539	101.74183	B1	16	0.00000
981	98 HIS	HD1	-46.72307	-40.60043	101.51181	Bl	16	0.00000 0.00000
982	98 HIS	CD2	-44.21286		102.89430	BI	16 16	0.00000
983	98 HIS	NE2	-44.02278	-38.92334	101.58753	D1	16	0.00000
984	98 HIS	CE1	-45.05268	-39.48082	106.05217		16	0.00000
985	98 HIS	C	-47.56750	-40.40062 -40.99638	106.99225	Bl	16	0.00000
986	98 HIS	0	-47.04279 -48.87987		105.81218	91	17	0.00000
987	99 PHE 99 PHE	И	-49.25243		105.00666	Bl	17	0.00000
. 988 989	99 PHE	H CA	-49.78029			Bl	17	0.00000
990	99 PHE	CB	-50.48946		107.59234	Bl	17	0.00000
991	99 PHE	CG	-49.63868	_ 39.72312	108.81497	Bl	17	0.00000
992	99 PHE	CD1	-48.50685	-38.87580	108.76543	B1	17	0.00000 0.00000
993	99 PHE	CD2		-40.47457	109.97439	B1	17 17	0.00000
994	99 PHE	CE1	-47.64103	-38.80002	109.87537		17	0.00000
995	99 PHE	CE2	-49.06252	-40.39779 -39.56658	111.02626	Bì	17	0.00000
996	99 PHE	CZ	-47.92137	-41.97404	106.19800	Bl	17	0.00000
997 998	99 PHE 99 PHE	ë	-51 48007	-41.71068	105.22101	Bl	17	0.00000
999	100 PRE	И О	-50.86837	-43.12653	200.000	Bl	18	0.00000
1000	100 PHE	·H	-50 32768	-43.27017	107.68944	Bl	18	0.00000
1001	100 PHÉ	CA	-51 84718	-44.10783	106.41132	Bl	18	0.00000 0.00000
1002	100 PHE	CB	-51.44468	-45.52210	106.84262	P.1	18 18	0.00000
1003	100 PHE	CG	-51.08740	-46.34931	105.63001	BI	18	0.00000
1004	100 PHE	CD1	-49.81159	-46.95065 -46.50158	103.55500	Bl	18	0.00000
1005	100 PHE	CD2	-52.01269	-47.69723	104.41336	B1	18	0.0000
1006 1007	100 PHE	CE1 CE2	-49.45011	-47.24829	103.42719	Bl	18	0.00000
1007	100 PHE	CZ	-50 37408	-47.83859	103.35608	Вl	18	0.00000
1009	100 PKE	c	-53 21727	-43.84401	106.96975	BI	18	0.00000
1010	100 PHE	0	-53 38235	-43.49692	108.13319	BI	18	0.00000 0.00000
1011	101 ASK	N	-54 19611	-44.08275	106.08672	3:	19 19	0.00000
1012	101 ASN	н	-53.92777	-44.28429 -43.97453	105 37772	ום	19	0.00000
1013	101 ASN	CA	-55.63451	-45.9/403 -45.03350	105.37273	P.1	19	0.00000
1014	101 ASN	CB	-56.35400	-45.03359 -44.83624	105.52911	21	19	0.00000
1015	101 ASN	CG	-51.00040	-44.06601	104.77523	21	19	0.0000
1016	101 ASN 101 ASN	OD1 ND2	_58 53 327	-45.52184	206.41339	9.1	19	0.00000
1017 1018	101 ASN	HD21	_58 04843	-46.21753	107.02623	= -	19	0.00000
1018	101 ASK	HD22	_50 50 666	-45.49785	106.46862	2.7	19	0.00000 0.00000
1020	101 AS:	С	-55.06277	-44.09262	107.83398	51	1 9	0.0000

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1021	101 A	SN O	16.7373	7 -43.2399	7 108.39463	B1	19	0.00000
1022					1 108.46454		20	0.00000
2025	102 G		-55.0037	0 -45.8218	8 108.01935	Bl	20	0.00000
1024	102 G		-55.9743	8 -45.3235	3 109.88157	Bl	20	0.00000
1025	102 G	LY C		8 -44.5987			20	0.00000
1026	102 G	LY O	-54.4600				20	0.00000
1027	103 T	HR N	-54.8474	4 -43.2971	6 110.54648	Bl	21	0.00000
1028	103 T	HR H	-55.3551	2 -42.9073	1 109.77435		21	0.00000
1029	103 T	HR CA		9 -42.3954			21	0.00000
1030	103 T	HR CB		7 -41.8126			21	0.00000
1031	103 T	HŘ OG1		5 -40.5335			21	0.00000
1032		HR HG1	-54.8197	4 -40.0991	9 113.49221	Bl	21	0.00000
1033		HR CG2	-54.8692	5 -42.7243	113.73959	B1	21	0.00000
1034	103 T	,-	-52.56983	3 - 42.92386	111.66956	Bl	21	0.00000
1035	103 T	_	-51.96086	-42.5928	112.68707	B1	21	0.00000
1036	104 GI		-52.05837	7 -43.7843	110.78327	D.7	22 22	0.00000
1037	104 G1		-52.50836	-43.9402	109.90068	10.1 10.1	22	0.00000
1038	104 GI 104 GI			7 -45.95191			22	0.00000
1039			-50.88647 51.13368				22	0.00000
1041	104 GI				113.43075		22	0.00000
1042	104 GI				114.57130		22	0.00000
1043	104 GI			-46.13148			22	0.00000
1044	104 GI			-43.91901			22	0.00000
1045	194 GL			-43.40196			22	0.00000
1046	105 AR			-44.04578			23	0.00000
1047	105 AR	•			111.97754	Bl	23	0.00000
1048	105 AR	. •			110.79677		23	0.00000
1049	105 AR			-43.14957	112.09184	Bl	23	0.00000
1050	105 AP	•		-42.37595		Bl	23	C.00000
1051	105 AR	•	-43.89570	-43.22016	111.65081	Bl	23	0.00000
1052	105 AR	G NE	-42.65640	-42.48704	111.89499	Bl	23	0.0000
1053	105 AR			-42.45915			23	0.00000
1054	105 AR		-42.05554	-41.77685	110.93892		23	0.00000
1055	105 AR			-41.00938			23	0.00000
1056	105 AR	•		-40.49136		Bl	23	0.00000
1057	105 AR		-40.73208	-40.92747	112.21979		23	0.00000 0.00000
1058	105 AR		-42.47661	-41.82718	109.68286		23 23	0.00000
1059	105 AR				109.00310	D)	23	0.00000
1060	105 AR			-42.57763	109.74376		23	0.00000
1061 1062	105 AR		-46.47373	-45 25363	109.97034	Bl	23	0.00000
1063	106 VA	.*	-46 32883	-43.23503	108.56827	Bl	24	0.00000
1064	106 VA	,			108.37187		24	0.00000
1065	106 VA				107.58352		24	0.00000
1066	106 VA		-46.27081	-44.39073	106.24298	B1	24	0.00000
1067	106 VA		-45.79579	-45.65575	105.52366	B1	24	0.00000
1068	706 AVI	CG2	-47.77990	-44.41166	106.42922	B1	24	0.00000
1069	106 AVI		-44.14065	-43.66075	107.41554	B1	24	0.00000
1070	106 AY		-43.30354	-43.80891	108.29696	Bl	24	0.00000
1071	107 ARC	•			106.29416		25	0.00000
1072	107 ARC		-44.58452	-42.69040	105.65671	BI	25	0.00000
1073	107 ARC		-42.49561	-42.53115	106.12329	BI	25	0.00000
1074	107 ARG		-41.95685	-43.01317	104.76032	81	25	0.00000 0.00000
1075	107 ARG		-41.96328	-42.02619	103.58362	B1	25	0.00000
1076	107 ARG		-42.33357	-42.66829	102.24883	D.I.	25 25	0.00000
1077	107 ARG		-43.76838	-42.93515	102.23147	BJ BJ	25	0.00000
1078	107 ARG		-44.3/946.	-42.14314	102.28284	B1	25	0.00000
1079 1080	107 ARG		-45.56598	-44 32747	102.35905	B 1	25	0.0000
1081	107 ARG	hhill	-46.08629	-45.19207	102.34562	5 1	25	0.00000
1082	107 ARG	HH12	-45.19539	-43.55730	102.54612	E I	25	0.00000
1093	107 ARG	11:12	-43.47983	-45.23695	102.05793	5 1	25.	0.00000
1084	107 ARG	8821	-43.86541	-45.15896	102.05452	21	25	0.00000

1085 107 ARG HH22 2, 49.06 -45.11987 101.9496 B1 25 0.00000 1086 107 ARG C	_ندعر.	೬೯೩೭ . ೧೯೨		೭೫# ೬¢೨	25 14:58:4	8 1993		79	
1086 107 ARG C	1085	107 APC	HH22	2,49404	-45.11997	101.94496	Bl	25	0.00000
1068 108 LEU N -40.12633 -40.58602 106.38627 B1 25 0.00000 1089 108 LEU C H -40.32617 -41.17985 106.29320 B1 26 0.00000 1091 108 LEU C B -40.80087 -38.80934 108.01248 B1 26 0.00000 1092 108 LEU C B -40.80087 -38.80934 108.01248 B1 26 0.00000 1093 108 LEU C B -40.80087 -38.80934 108.01248 B1 26 0.00000 1093 108 LEU C D -41.50347 -37.89689 108.38574 B1 26 0.00000 1093 108 LEU C D1 -42.75820 -38.44729 109.53808 B1 26 0.00000 1095 108 LEU C D2 -41.50070 -36.44650 108.50161 B1 26 0.00000 1095 108 LEU C D3 -41.50070 -36.44650 108.50161 B1 26 0.00000 1095 108 LEU C -39.62061 -38.77273 105.85764 B1 26 0.00000 1097 109 LEU N -39.52150 -37.47828 105.58773 B1 27 0.00000 1099 109 LEU N -39.52150 -37.47828 105.58773 B1 27 0.00000 1099 109 LEU N -39.52150 -37.47828 105.58773 B1 27 0.00000 1099 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.30683 -37.02679 104.90334 B1 27 0.00000 1100 109 LEU C -38.8411 -34.77322 105.47943 B1 27 0.00000 1100 109 LEU C -38.88411 -34.77322 105.47943 B1 27 0.00000 1100 109 LEU C -37.59969 -35.58061-105.21266 B1 27 0.00000 1100 100 LEU C -37.59969 -35.58061-105.21266 B1 27 0.00000 1100 109 LEU C -37.59969 -35.58061-105.21266 B1 27 0.00000 1100 1100 LEU C -38.88411 -34.77322 105.47943 B1 27 0.00000 1100 1100 LEU C -38.8				2.37787	-41.03546	106.32504	Bl		
1088 108 LEU N				-43.36896	-40.31981	106.43645	Bl		
1090 108 LEU CA -40,89942 -39,15744 106,53751 B1 26 0,00000 1091 108 LEU CG -41,96347 -37,89659 108,35574 B1 26 0,00000 1093 108 LEU CD2 -41,5820 -38,474729 109,53808 B1 26 0,00000 1095 108 LEU CD2 -41,5820 -38,474729 109,53808 B1 26 0,00000 1096 108 LEU CD2 -39,62061 -38,77373 105,58764 B1 26 0,00000 1097 109 LEU N -39,52150 -37,47828 105,56773 B1 27 0,00000 1098 109 LEU C -38,36683 -37,02679 104,90334 B1 27 0,00000 1099 109 LEU CA -38,36683 -37,02679 104,90334 B1 27 0,00000 1000 109 LEU CG -39,74480 -36,55655 102,82634 B1 27 0,00000 1001 109 LEU CG -39,74490 -36,55655 102,82634 B1 27 0,00000 1002 109 LEU CD1 -39,42490 -36,55655 102,82634 B1 27 0,00000 1003 109 LEU CD2 -40,49190 -37,56523 101,91138 B1 27 0,00000 1004 109 LEU CD3 -39,74490 -36,55655 102,82634 B1 27 0,00000 1105 109 LEU CD4 -38,88411 -34,77322 105,47943 B1 27 0,00000 1106 100 LEU CD5 -37,9959 -35,56651 105,21268 B1 27 0,00000 1107 110 GLU M -36,01498 -36,01766 105,00992 B1 28 0,00000 1109 110 GLU CD5 -33,83808 -31,8716 106,30413 B1 28 0,00000 1109 110 GLU CD5 -33,37957 -30,74951 106,60993 B1 28 0,00000 1111 110 GLU CD5 -33,37957 -30,74951 106,60993 B1 28 0,00000 1112 110 GLU CD5 -33,37957 -33,47627 106,49578 B1 28 0,00000 1113 110 GLU CD5 -33,37957 -34,7895 106,36491 B1 28 0,00000 1114 110 GLU CD5 -33,37957 -34,7895 106,36491 B1 28 0,00000 1112 111 GLU CD5 -33,37957 -34,7895 106,36491 B1 28 0,00000 1113 110 GLU CD5 -33,37957 -34,7895 0,00000 -34,28577 -34,8968 -34,8977 -33,4975 -34,8975 -34,8977 -34,8977 -34,8977 -34,		108 LEU	N	-41.12633	-40.58602	106.39627	B1		
1091 108 LEU CB	1089			- 40.32617	-41.17985	106.29320	81 B1		•
1092 108 LEU CG				-40.89942	-39.15/44				
1093 108 LEU CD1 -42,75820 - 38, 44729 109,53808 B1 26 0.00000 1094 108 LEU CD -41,55070 - 36, 44650 108, 50151 B1 26 0.00000 1095 108 LEU CD -38,62061 - 38,77273 105,85764 B1 26 0.00000 1097 109 LEU N -39,52150 - 37,47828 105,56773 B1 27 0.00000 1099 109 LEU CA -40,28459 - 36,84167 105,72467 B1 27 0.00000 1099 109 LEU CA -38,30683 - 37,02679 104,90334 B1 27 0.00000 1010 109 LEU CB -38,47418 -37,24039 103,38009 B1 27 0.00000 1010 109 LEU CD2 -39,47449 -36,55951 102,86736 B1 27 0.00000 1010 109 LEU CD2 -40,4490 -36,55951 102,86736 B1 27 0.00000 103 109 LEU CD2 -40,4490 -36,55651 102,17066 B1 27 0.00000 105 109 LEU CD2 -40,49190 -37,56523 101,91138 B1 27 0.00000 105 109 LEU CD2 -40,49190 -37,56523 101,91138 B1 27 0.00000 105 109 LEU CD -37,99969 -35,58061-105,21268 B1 27 0.00000 105 109 LEU CD -36,70092 -35,29768 105,12126 B1 20 0.00000 107 110 GLU B -36,01488 -36,01766 105,0092 B1 28 0.00000 1101 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 1112 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 1113 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 1114 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 1115 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 1116 111 ARG R -35,580977 -32,468047 101,20352 B1 28 0.00000 1116 111 ARG R -35,580977 -32,468047 101,20352 B1 28 0.00000 1115 110 GLU CD -33,83808 -31,87178 106,30413 B1 28 0.00000 112 111 ARG R -35,580977 -33,493076 104,115500 B1 28 0.00000 112 111 ARG R -35,580977 -33,493076 104,115500 B1 28 0.00000		•		-40.80087	-30.80934	108.35874	B1		
1095 108 LEU CD2 -41.50070 -36.44650 108.50151 B1 26 0.00000 1096 108 LEU CD -39.62061 -38.77273 105.85764 B1 26 0.00000 1097 109 LEU N -39.52150 -37.47428 105.56773 B1 27 0.00000 1098 109 LEU H -40.28459 -36.84167 105.72467 B1 27 0.00000 1009 109 LEU CB -38.30683 -37.02679 104.90334 B1 27 0.00000 1010 109 LEU CB -38.30683 -37.02679 104.90334 B1 27 0.00000 1010 109 LEU CG -39.74498 -36.55695 102.82634 B1 27 0.00000 102 109 LEU CG -39.74490 -36.55695 102.82634 B1 27 0.00000 103 109 LEU CC -37.99969 -35.56523 102.17066 B1 27 0.00000 104 109 LEU CC -37.99969 -35.56523 102.17066 B1 27 0.00000 1105 109 LEU C -37.99969 -35.56523 102.82634 B1 27 0.00000 1106 110 GLU N -36.70092 -35.756523 101.91138 B1 27 0.00000 1107 110 GLU N -36.70092 -35.75661 -35.12126 B1 28 0.00000 1108 110 GLU CD -35.50077 -33.47627 106.49578 B1 28 0.00000 1109 110 GLU CD -33.83088 -31.87178 106.49578 B1 28 0.00000 1111 110 GLU CD -33.83088 -31.8718 106.30413 B1 28 0.00000 1112 110 GLU CD -33.83088 -31.8718 106.10521 B1 28 0.00000 1114 110 GLU CD -33.83088 -31.8718 106.10521 B1 28 0.00000 1115 110 GLU CD -33.83088 -31.8718 106.10521 B1 28 0.00000 1116 111 ARG R -35.61365 -32.16766 103.37387 B1 28 0.00000 1116 111 ARG R -35.61365 -32.16766 103.37387 B1 29 0.00000 1116 111 ARG R -35.61365 -32.16766 103.37387 B1 29 0.00000 1120 111 ARG G -35.61365 -32.16766 103.37387 B1 29 0.00000 1121 111 ARG R -35.61365 -32.16766 103.37387 B1 29 0.00000 1122 111 ARG R -37.71660 -37.8683 -38.4683 -38.4683 -38.4683 -38.4683 -				-42.75820	-38.44729	109.53808	B1	26	0.00000
1095 108 LEU C -39,62061 -38,77273 105,85764 B1 26 0.00000 1097 109 LEU N -39,52150 -37,47828 105,56773 B1 27 0.00000 1098 109 LEU N -40,28459 -36,84167 105,72467 B1 27 0.00000 1100 109 LEU CA -38,30683 -37,02679 104,90334 B1 27 0.00000 1101 109 LEU CA -38,47418 -37,24039 103,39009 B1 27 0.00000 1102 109 LEU CA -39,47490 -36,55895 102,87634 B1 27 0.00000 1102 109 LEU CA -39,42210 -35,55831 102,17066 B1 27 0.00000 1103 109 LEU CD -40,49190 -37,55523 101,91138 B1 27 0.00000 1104 109-LEU C -37,99969 -35,55623 101,91138 B1 27 0.00000 1105 109 LEU C -37,99969 -35,55623 101,91138 B1 27 0.00000 1105 109 LEU C -37,99969 -35,55623 101,91138 B1 27 0.00000 1105 109 LEU C -37,99969 -35,55623 101,91138 B1 27 0.00000 1105 101 GLU II -36,-01498 -36,01786 105,00992 B1 28 0.00000 1106 110 GLU II -36,-01498 -36,01786 105,00992 B1 28 0.00000 1109 110 GLU CA -36,20315 -33,32985 105,08184 B1 28 0.00000 1101 110 GLU CA -35,32216 -32,02210 106,60993 B1 28 0.00000 1112 110 GLU CA -33,83808 -31,87178 106,60933 B1 28 0.00000 1113 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -33,83808 -31,87178 106,10521 B1 28 0.00000 1114 110 GLU CA -35,6565 -32,16666 -39,				-41.50070	-36.44650	108.50151	Bl	26	
1097 109 LEU N -39.52150 -37.47828 105.56773 B1 27 0.00000 109 LEU H -40.28559 -36.84167 105.72467 B1 27 0.00000 109 LEU CB -38.30683 -37.26279 104.90334 B1 27 0.00000 1101 109 LEU CB -38.47418 -37.24039 103.39009 B1 27 0.00000 1102 109 LEU CD -39.74490 -36.55965 102.82634 B1 27 0.00000 1103 109 LEU CD -39.74490 -36.55965 102.82634 B1 27 0.00000 1103 109 LEU CD -39.74490 -36.55965 102.87634 B1 27 0.00000 1103 109 LEU CD -39.74490 -36.55965 102.87634 B1 27 0.00000 1104 109-LEU CD -39.4490 -37.55523 101.91138 B1 27 0.00000 1104 109-LEU CD -38.88411 -34.77322 105.47943 B1 27 0.00000 1106 110 GLU N -36.01498 -36.01498 -36.5664 -305.12126 B1 28 0.00000 1106 110 GLU CD -36.88411 -34.77322 105.47943 B1 28 0.00000 1108 110 GLU CD -35.80977 -33.47627 106.6993 B1 28 0.00000 1101 110 GLU CD -35.80977 -33.47627 106.6993 B1 28 0.00000 1113 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1113 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1114 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1114 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1114 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1114 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.83088 -31.87178 106.30413 B1 28 0.00000 1115 110 GLU CD -35.65651 -35.65651 -35.6667				-39.62061	-38.77273	105.85764	Bl		
1098 109 LEU N				-38.76368	-39.61093	105.60154	Bl		
109 109 LEU CA -38, 30683 -37, 02679 104, 90334 B1 27 0.00000	1097	109 LEU		-39.52150	-37.47828	105.56773	B1		
1100 109 LEU CB -38. 47418 -37. 24039 103.39009 B1 27 0.00000 1101 109 LEU CD -39. 44290 -36. 55995 102. 82634 B1 27 0.00000 1103 109 LEU CD -39. 42210 -35. 25631 102. 17066 B1 27 0.00000 1104 109 LEU CD -40. 49190 -37. 56523 101. 91138 B1 27 0.00000 1105 109 LEU C -37. 99969 -35. 58061 105. 21268 B1 27 0.00000 1106 110 GLU N -36. 70092 -35. 358061 105. 21268 B1 27 0.00000 1107 110 GLU N -36. 70092 -35. 358061 105. 09192 B1 28 0.00000 1108 110 GLU C -36. 80411 -34. 77322 105. 7943 B1 28 0.00000 1109 110 GLU C -36. 80418 -36. 01786 105. 00192 B1 28 0.00000 1109 110 GLU C -35. 80977 -33. 47627 106. 49578 B1 28 0.00000 1111 110 GLU C -33. 33216 -32. 02210 106. 60993 B1 28 0.00000 1112 110 GLU C -33. 337957 -30. 74951 106. 10521 B1 28 0.00000 1113 110 GLU C -35. 01026 -33. 94076 104. 15500 B1 28 0.00000 1115 111 GGU C -34. 25577 -34. 90807 104. 11555 B1 28 0.00000 1115 111 ARG R -35. 61365 -32. 16766 103. 37327 B1 29 0.00000 1117 111 ARG R -35. 61365 -32. 16766 103. 37327 B1 29 0.00000 1119 111 ARG CB -33. 99004 -33. 60407 101. 20352 B1 29 0.00000 1120 111 ARG CB -33. 99004 -33. 60407 101. 20352 B1 29 0.00000 1121 111 ARG CB -33. 99004 -33. 60407 101. 20352 B1 29 0.00000 1122 111 ARG H -35. 61365 -32. 16766 103. 37327 B1 29 0.00000 1123 111 ARG H -36. 61365 -32. 16766 103. 37327 B1 29 0.00000 1124 111 ARG CB -35. 61366 -35. 70037 100. 25355 B1 29 0.00000 1125 111 ARG H -36. 61365 -36. 68971 99. 97. 97. 97. 97. 97. 97. 97. 97. 97.				-40.28459	-36.84167	105.72467	BJ DI		
1101 109 LEU CD				-38.30683	-37.02679	103.30009	Bl		
1102				-30.5/418	-36.59695	102.82634	Bl		
1103 109 LEU CD2				-39.42210	-35,25631	102.17066	Bl	27	0.00000
104 109 LEU	_			-40.49190	-37.56523	101.91138	Bl		
1105 109 LEU O -38.88411 -34.77322 105.47943 B1 27 0.00000 1106 110 GLU N -36.70092 -35.29768 105.12126 B1 28 0.00000 1107 110 GLU CA -36.20315 -33.92985 105.00929 B1 28 0.00000 1109 110 GLU CC -35.80977 -33.47627 106.49578 B1 28 0.00000 1110 110 GLU CC -35.32216 -32.02210 106.60993 B1 28 0.00000 1111 110 GLU CC -35.33216 -32.02210 106.60993 B1 28 0.00000 1112 110 GLU CD -33.83808 -31.87178 106.30413 B1 28 0.00000 1113 110 GLU CD -33.83808 -31.87178 106.30413 B1 28 0.00000 1113 110 GLU CD -33.37957 -30.74951 106.10521 B1 28 0.00000 1114 110 GLU C -35.01026 -33.94076 104.15500 B1 28 0.00000 1115 110 GLU C -34.25577 -34.90807 104.15500 B1 28 0.00000 1116 111 ARG N -34.89439 -32.86843 103.37222 B1 29 0.00000 1117 111 ARG H -35.61365 -32.16766 103.37387 B1 29 0.00000 1118 111 ARG CA -33.75645 -32.74903 102.46249 B1 29 0.00000 1119 111 ARG CC -35.45106 -33.70037 100.75305 B1 29 0.00000 1120 111 ARG NC -35.67880 -34.83036 99.75174 B1 29 0.00000 1121 111 ARG NC -35.67880 -34.83036 99.75174 B1 29 0.00000 1122 111 ARG NC -35.75098 -34.61037 100.27575 B1 29 0.00000 1123 111 ARG NR -37.70984 -35.61357 100.22757 B1 29 0.00000 1125 111 ARG NR -37.70986 -34.61357 100.22757 B1 29 0.00000 1125 111 ARG NR -37.75098 -36.3238* 99.09789 B1 29 0.00000 1126 111 ARG NR -37.75098 -36.3238* 99.09789 B1 29 0.00000 1127 111 ARG HH11 -39.12269 -37.49888 98.68276 B1 29 0.00000 1128 111 ARG HH11 -39.43712 -36.16699 99.70940 B1 29 0.00000 1128 111 ARG HH21 -36.63931 -37.06889 99.79189 B1 29 0.00000 1133 112 CYS N				-37.99969	-35.58061·	·105.21268	81		
1107 110 GLU II		109 LEU	0	-38.88411	-34.77322	105.47943	B1		
1108	1106			-36.70092	-35.29768	105.12126	BI		
110				-36.01498	-36.01786	105.00992	BJ DT		
1110				-36.20315	-33.92903	105.00104	B1		
1111 110 GLU CD				-35.809//	-32 02210	106.60993	81		0.00000
1112 110 GLU 0E1				-33 B380B	-31.87178	106.30413	вı	28	0.00000
1113 110 GLU OE2 -33 09671 -32.85201 106.36491 B1 28 0.00000 1114 110 GLU C -35 01026 -33.94076 104.11580 B1 28 0.00000 1115 110 GLU O -34.25577 -34.90807 104.11585 B1 28 0.00000 1116 111 ARG N -34.89439 -32.86843 103.37222 B1 29 0.00000 1117 111 ARG H -35.61365 -32.16766 103.37387 B1 29 0.00000 1118 111 ARG CA -33.75665 -32.747903 102.46249 B1 29 0.00000 1119 111 ARG CB -33.99004 -33.60407 101.20352 B1 29 0.00000 1120 111 ARG CB -35.45106 -35.70037 100.75305 B1 29 0.00000 1121 111 ARG CD -35.67880 -34.83036 99.75174 B1 29 0.00000 1122 111 ARG NE -37.09124 -35.21147 99.72614 B1 29 0.00000 1123 111 ARG NE -37.71860 -34.61357 100.22757 B1 29 0.00000 1124 111 ARG CC -37.50098 -36.32384 99.09789 B1 29 0.00000 1125 111 ARG NH11 -39.12269 -37.49888 98.68276 B1 29 0.00000 1126 111 ARG NH11 -39.43712 -36.668971 99.16119 B1 29 0.00000 1127 111 ARG NH2 -36.63391 -37.06893 98.41518 B1 29 0.00000 1128 111 ARG NH2 -36.63391 -37.06893 98.41518 B1 29 0.00000 1129 111 ARG HH22 -35.67503 -36.78688 98.36264 B1 29 0.00000 1130 121 ARG HH22 -35.67503 -36.78688 98.36264 B1 29 0.00000 1131 111 ARG C -33.49135 -31.31194 102.08170 B1 29 0.00000 1133 112 CYS N -32.20568 -30.97057 102.12414 B1 30 0.00000 1135 112 CYS CA -31.80468 -29.62360 101.73826 B1 30 0.00000 1136 112 CYS CA -31.80468 -29.62360 101.73826 B1 30 0.00000 1137 112 CYS CA -31.80468 -29.62365 102.91930 B1 30 0.00000 1138 112 CYS CA -31.80468 -29.62360 101.73826 B1 30 0.00000 1139 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.000000 1131 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.000000 1131 112 CYS C -				-33.37957	-30.74951			28	
1114 110 GLU C				-33.09671	-32.85201	106.36491	Bl		
1115 110 GLU O				-35.01026	-33.94076	104.15500	Bl		
1116				-34.25577	-34.90807	104.11585	B1		
1118 111 ARG CA	1116	111 ARG	Й	-34.89439	-32.86843	103.37222	B.T		
1118 111 ARG CA	1117			-35.61365		103.3/38/	B1		
1120 111 ARG CG				-33.75645	-32.74903	101 20352	B1		
1121 111 ARG CD			•	-35.99004	-33.70037	100.75305	Bl		0.00000
1122 111 ARG				-35.67880	-34.83036	99.75174	Bl	29	
1123				-37.09124	-35.21147	.99.72614	B1		
1125 111 ARG				-37.71860	-34.61357	100.22757	B1		
1126 111 ARG HH11 -39.12269 -37.49888 98.68276 B1 29 0.00000 1127 111 ARG HH12 -39.43712 -36.16699 99.70940 B1 29 0.00000 1128 111 ARG NH2 -36.63391 -37.06893 98.41518 B1 29 0.00000 1129 111 ARG HH21 -36.91501 -37.90348 97.94325 B1 29 0.00000 1130 111 ARG HH22 -35.67503 -36.78688 98.36264 B1 29 0.00000 1131 111 ARG C -33.49135 -31.31194 102.08170 B1 29 0.00000 1132 111 ARG C -33.49135 -31.31194 102.08170 B1 29 0.00000 1133 112 CYS N -32.20568 -30.97057 102.12414 B1 30 0.00000 1134 112 CYS H -31.50733 -31.64380 102.37455 B1 30 0.00000 1134 112 CYS CA -31.80468 -29.62360 101.73826 B1 30 0.00000 1136 112 CYS CB -31.12874 -28.92365 102.91930 B1 30 0.00000 1137 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1138 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1139 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1140 113 ILE N -31.15975 -28.70696 99.65078 B1 31 0.00000 1140 113 ILE N -31.84193 -28.00116 99.86111 B1 31 0.00000 1141 113 ILE C -33.55306 -28.70228 98.32464 B1 31 0.00000 1144 113 ILE C -33.55306 -28.70228 98.32464 B1 31 0.00000 1144 113 ILE C -33.55306 -28.70228 98.32464 B1 31 0.00000 1144 113 ILE C -33.55306 -28.70228 98.32464 B1 31 0.00000 1144 113 ILE C -33.00539 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE C -33.00539 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE C -33.00539 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE C -33.00539 -29.37434 97.82323 B1 31 0.00000 1146 113 ILE C -33.44047 -29.50025 94.93272 91 31 0.00000 1147 113 ILE C -33.44047 -29.50025 94.93272 91 31 0.00000 1147 113 ILE C -33.44047 -29.50025 94.93272 91 31 0.00000 1147 113 ILE C -33.44047 -29.50025 94.93272 91 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31 0.00000 1147 113 ILE C -33.44047 -29.50025 94.93272 91 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31 0.00000 1147 113 ILE C -33.00576 -27.29665 97.96536 B1 31	1124	111 ARG	CZ	-37.50098	-36.32384	99.09789	B1		
1127 111 ARG				-38.78116	-36.68971	99.10119	R1		
1128 111 ARG NH2				-39,12269	-37.49888				
1120				-35.43/12	-37 06893				0.00000
1130 111 ARG		: -	NU21	-36.03331	-37.90348		Bl	29	0.00000
1131 111 ARG C			HH22	-35.67503	-36.78688	.98.36264	B1	29	
1132 111 ARG O				-33.49135	-31.31194	102.08170	B1		
1133 112 CYS H			0	-34.39395	-30.54569	101.76699	Bl		
1134 112 CYS CA -31.80468 -29.62360 101.73826 B1 30 0.00000 1136 112 CYS CB -31.12874 -28.92365 102.91930 B1 30 0.00000 1137 112 CYS SG -30.70297 -27.19555 102.57697 B1 30 0.00000 1138 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1138 112 CYS C -30.87388 -29.64998 100.54380 B1 30 0.00000 1139 112 CYS O -29.97769 -30.48052 100.40406 B1 30 0.00000 1140 113 ILE N -31.15975 -28.70696 99.65078 B1 31 0.00000 1141 113 ILE H -31.84193 -28.00116 99.86111 B1 31 0.00000 1142 113 ILE CA -30.55306 -28.70228 98.32464 B1 31 0.00000 1143 113 ILE CB -31.56021 -29.33526 97.30706 B1 31 0.00000 1144 113 ILE CB -33.00339 -29.37434 97.82323 B1 31 0.00000 1144 113 ILE CG2 -33.00339 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE CG1 -31.52431 -28.74184 95.89635 B1 31 0.00000 1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE CD -33.08576 -27.29665 97.96536 B1 31 0.00000	1133	112 CYS	N	-32.20568	-30.97057	102.12414	B1		
1135 112 CYS CA		•	•	-31.50733	-31.64380	102.3/455	BI		
1137 112 CYS SG				-31.80468	-29.02360	102.73820	B1		
1138 112 CYS C				-31.12077	-27.19555	102.57697	Bl		0.00000
1139 112 CYS O				-30.87388	-29.64998	100.54380	Bl	30	
1140 113 ILE N -31.15975 -28.70696 99.65078 B1 31 0.00000 1141 113 ILE H -31.84193 -28.00116 99.86111 B1 31 0.00000 1142 113 ILE CA -30.55306 -28.70228 98.32464 B1 31 0.00000 1143 113 ILE CB -31.56021 -29.33526 97.30706 B1 31 0.00000 1144 113 ILE CG2 -33.00339 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE CG1 -31.52431 -28.74184 95.89635 B1 31 0.00000 1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE CD -33.08576 -27.29665 97.96536 B1 31 0.00000				-29.97769	-30.48052	100.40405	Bl		
1141 113 ILE H -31.84193 -28.00116 99.86111 B1 31 0.00000 1142 113 ILE CA -30.55306 -28.70228 98.32464 B1 31 0.00000 1143 113 ILE CB -31.56021 -29.33526 97.30706 B1 31 0.00000 1144 113 ILE CG2 -33.00339 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE CG1 -31.52431 -28.74184 95.89635 B1 31 0.00000 1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE C -30.08576 -27.29665 97.96536 B1 31 0.00000				-31.15975	-28.70696	99.65078	B1		
1142 113 ILE CB -31.56021 -29.33526 97.30706 B1 31 0.00000 1144 113 ILE CG2 -33.00319 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE CG1 -31.52431 -26.74184 95.89635 B1 31 0.00000 1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE C -30.08576 -27.29665 97.96536 B1 31 0.00000		113 ILE	H	-31.84193	-28.00116	99.86111	51		
1143 113 ILE CG2 -33.00319 -29.37434 97.82323 B1 31 0.00000 1145 113 ILE CG1 -31.52431 -26.74184 95.89635 B1 31 0.00000 1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE C -30.08576 -27.29665 97.96536 B1 31 0.00000				-30.55306	-28.70228	70.32404 70706	BJ DI		
1145 113 TLE CG1 -31.52431 -26.74184 95.89635 B1 31 0.00000 1146 113 TLE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 TLE C -30.08576 -27.29665 97.96536 B1 31 0.00000				-31.56021	-29.33526				
1146 113 ILE CD -32.44047 -29.50025 94.93272 B1 31 0.00000 1147 113 ILE C -30.08576 -27.29665 97.96536 B1 31 0.00000				-33.00329	-28.74184				0.00000
1147 113 ILE C -30.08576 -27.29665 97.96536 B1 31 0.00000		•		-32.44047	-29.50025	94.93272	81	31	
				-30.08576	-27.29665	97.96536	31		
				-30.75333	-26.29900	98.21317	Bl	31	0.00000

./DR1	_xr:	7. CZ		ಮಿಶ್	حمة	25 14:	58:	8 1	993		19	
1149	. 11	4 7YF		20	8701	0 -27.2			7.4166	.	32	0.00000
1150		4 TY				0 - 27.2 0 - 28.1		, ,	7.4166.	וחינ	32	0.00000
1251		4 TYP				6 ~25.9			7.1025 7.1146		32	0.00000
1152		4 TYF				B -26.0			7.7134		32	0.00000
1153		4 777				7 -24.7			7.6131		32	0.00000
1154		4 7YF				-23.4			7.7906		32	0.00000
1155		4 TYP				5 -22.2			7.7526		32	0.00000
1156		4 7YR				-24.8			7.3875		32	0.00000
1157		4 TYP				-23.6			7.3458		32	0.00000
1158	11					-22.3			7.53154		32	0.00000
1159		4 TYR				-21.2			7.5097		32	0.00000
1160	11					-21.4			7.21984		32	0.00000
1161		4 TYR				-25.8			5.61430		32	0.00000
1162	11					-26.7			1.91236		32	0.00000
1163		5 ASN	-			-24.6			12919		33	0.00000
1164		5 ASN				-24.0			77643		33	0.00000
1165		5 ASN				-24.4			.68566		33	0.00000
1166		5 ASN	CB			-23.8			.27078		33	0.00000
1167		5 ASN	CG			-22.43			.78440		33	0.00000
1168		S · ASN		26.5							33	0.00000
1169		5 ASN	ND2			-21.48			.96271		33	0.00000
1170		5 ASN		-28.0					.14797		33	0.00000
1171		NZA 6	HD22			-20.52			.13946		33	0.00000
1172		ASN	С			-25.59		92	.77321	Bl	33	0.00000
1173		ASN	ō			-25.91			.85250		33	0.00000
2174		GLN	N			-26.23			.07767		34	0.00000
1175	-	CIN	н			-25.92		93	.84830	Bl	34	0.00000
1176		GLN	CA			-27.38		92	.31212	Bl	34	0.00000
1177		GLN	CB			-26.98		90	.82631	B1	34	0.00000
1178		GLN	CG			-27.96		89	.80983	Bl	· 34	0.00000
1179	•	GLN	CD			-27.31		88	.43588	Bl	34	0.00000
1180	116	GLN	OEl	-32.3	2002	-27.33	002	87	.70605	Bl	34	0.00000
1181	116	GLM	NE2	-30.20	1428	-26.71	450		.08185		34	0.00000
1182	116	GLN	HE21	-29.3	9553	-26.71	.396		.67007		34	0.00000
1183		GIM	HE22	-30.1					.20205		34	0.00000
1184		GLN	С.			-28.73			.53043		34	0.00000
1185		GLN	٥			-29.77			.10767		34	0.00000
1186		GLU	N		•	-28.72			.24838		35	0.00000
1187	117	-	H			-27.88			.59525		35	0.00000
1188	117		CA			-30.02			.58636		35	0.00000
1189	117	· .	CB			-30.02			. 23572		35	0.00000
1190		CLU	CG			-29.71			.75724		35	0.00000
1191		GLU	CD			-29.87			.43636 .37359		35 35	0.00000
1192 1193		ern ern	OE1 OE2	•		-30.40 -29.45			.24135		35	0.00000
1194		CLU	C .			-30.39			.05360		35	0.00000
1195		GLU	0.	-28.39					. 93593		35	0.00000
1196	118		N			-31.70			31151		36	0.00000
1197		CLÜ	H	-28.37					57856		36	0.00000
1198	118		CA	-28.64				96.	70504	B1	36	0.00000
1199		GLU	CB	-28.86				96.	80220	Bl	36	0.00000
1200	11B	GLÜ	CG	-30.04				96	04454	Bl	36	0.00000
1201	228	GLŲ	CD	-30.18				96.	39969	Bl	36	0.00000
1202	118	GĮŲ	OE1	-31.31	366	-36.18	552	96.	62004	Bl	36	0.00000
1203		GLU	OE2	-29.17				96.	46236	81	36	0.00000
1204	118	CLU	C.	-27.45					59954		36	0.00000
1205		GLÜ		-26.30					22004		36	0.00000
1206	119	SER		-27.77					81671		37	0.00000
1207	119			-28.73					10568		37	0.00000
1208	119	SER		-26.67					74333		37	0.00000
1209	119			-26.79					25438		37	0.00000
1210	119	SER		-25.62					96854		37	0.00000
1211	119			-25.73					30354		37	0.00000
1212	119	SIR	С	-26.649	667	-32.108	169	100.	87999	91	37	0.00005

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1213	119 SER	٥	>5 72568	-32.90554	101 03211	81	37	0.00000
1213	120 VAL	и		-32.06808			38	0.00000
1215	120 VAL	н		-31.45938			38	0.00000
1216	120 VAL	ca ca		-32.99802			38	.0.00000
1217	120 VAL	СВ	-27.44790	-32.32244	104.13888	BJ	38	0.00000
1218	120 VAL	CG1	-25.93252	-32.18078	104.29388	Bl	38	0.00000
1219	120 VAL	CG2	-28.15631	-30.97900	104.33500	Bl	38	0.00000
1220	120 VAL	С	-29.23777	-33.53625	102.89372	Bl	38	0.00000
1221	120 VAL	0	-30.19812	-32.88656	102.49711	Bl	38	0.00000
1222	121 ARG	N	-29.34164	-34.75356	103.41944	Bl	39	0.00000
1223	121 ARG	н	-28.52525	-35.27675	103.67737	B1	39	0.00000
1224	121 ARG	CA	-30.65941	-35.37282	103.49447	BI	39	0.00000
1225	121 ARG	CB	-30.83879	-36.26556	102.25451	BI	39	0.00000 0.00000
1226	121 ARG	CG	-32.24891	-36.83907	102.13977	BI	39 39	0.00000
1227	121 ARG	CD	-32.36448	-38.11228	101.31436	B1	39	0.00000
122B	121 ARG	NE	-33,46636	-38.89614	101.86565	D)	39	0.00000
1229	121 ARG	HE	-33.58314	-38.82444	102.86363	20.1	39	0.00000
1230	121 ARG	C2	-34.21779	-39.71705	101.13235	B1	39	0.00000
1231	121 ARG	NH1	-35.18842	-40.40193	101.73310	B1	39	0.00000
1232	121 -ARG	HHII	-35.77971	-41.03441	102.23700	Bl	39	0.00000
1233	121 ARG		-33.33365	-39.84536	99.82444	BI	39	0.00000
1234	121 ARG	NH2			99.25675		39	0.00000
1235	121 ARG		-34.530 <u>7</u> 9	-30.37130	99.39731		39	0.00000
1236	121 ARG	HH22	-30.27034	-36.21013			39	0.00000
1237	121 ARG 121 ARG	C O	-20 86515	-36.82872	105.23483	Bl	39	0.00000
1238 1239	122 PHE	N	-32 04075	-36.24095	105.26404	Bl	40	0.00000
1240	122 PHE	н	-32.75061	-35.63810	104.89111	Bl	40	0.00000
1241	122 PHE	CA	-32.40668	-37.26064	106.24769	Bl	40	0.00000
1242	122 PHE	СВ	-33.75724	-36.87165	106.85296	Bl	40	0.00000
1243	122 PHE	CG	-33.64992	-36.46831	108.3041B	Bl	40	0.00000
1244	122 PHE	CDI	-32.56100	-35.69712	108.77664	Bl	40	0.00000
1245	122 PHE	CD2	-34.66985	-36.87929	109.19094	Вl	40	0.00000
1246	122 PHE	CE1	-32.49275	-35.34303	110.14035	Вl	40	0.00000
1247	122 PHE	CE2	-34,60187	-36.52254	110.55431	Bl	40	0.00000
1248	122 PHÈ	CZ	-33.51285	-35.75823	111.02379	Bl	40	0.00000
1249	122 PHE	С	-32.57844	-38.62424	105.60697	B1	40	0.00000
1250	122 PHE	0	-33.34168	-38.79208	104.65587	Bl	40	0.00000
1251	123 ASP	N	-31.06201	-39.60796	106.15389	B1	41 41	0.00000
1252	123 ASP	н	-31.23184	-39.47015	106.92346	B1	41	0.00000
1253	123 ASP	CA	-32.08552	-40.93522	105.36023	ים	41	0.00000
1254	123 ASP	CB	-30.831/1	-41.81726 -41.56041	104.59405	Bl	41	0.00000
1255	123 ASP	CC	-29.93161	-41.11310			41	0.00000
1256	123 ASP	OD1	-30.34905	-41 80191	103.45856	Bl	41	0.00000
1257 1258	123 ASP 123 ASP	C	-33.32362	-41 63618	106.09965	Bl	41	0.00000
1258	123 ASP	0	-34.00966	-41.21121	107.02319	Bl	41	0.00000
1260	124 SER	N	-33.62443	-42.74250	105.41429	B 1	42	0.00000
1261	124 5ER	н	-32.95329	-43.10971	104.76900	Bl	42	0.00000
1262	124 SER	CA	-34.94425	-43.35498	105.58880	Bl	42	0.00000
1263	124 SER	СВ	-35,18779	-44.37028	104.46149	Bl	42	0.00000
1264	124 SER	OG	-36.57945	-44.70776	104.37671	Bl	42	0.00000
1265	124 SER	HG	-36.91089	-44.87590	105.27413	Bl	42	0.00000 0.00000
1266	124 SER	С	-35.21640	-44.01057	106.93634	B1	42	0.00000
1267	124 SER	0	-36.33538	-44.42223	107.22372	Bl	42	0.00000
1268	125 ASP	N	-34.16447	-44.10325	107.74330	E.	43 43	0.00000
1269	125 ASP	Н	-33.26228	-43.73425	107.50494	b,	43	0.00000
1270	125 ASP	CA	-34.30492	-44.60689	109.104/1	נם	43	0.00000
1271	125 ASP	CB	-32.96210	-45.24580	109.50020	D 1	43	0.00000
1272	125 ASP	CG	-31.82155	13 10557	100 89075	91	43	0.00000
1273	125 ASP	021	-31.89194 -30.87249	-43.19333	110 29614	21	43	0.00000
1274	125 ASP	002	-30.87249 -34.69374	-47.430/0	110.30428	<u> </u>	43	0.00000
1275	125 ASP	C	-34.69374	-43.36644	111.22872	21	43	0.00000
1276	125 ASP	0	-35.10540	133471				

./DR1_	_೬೭೫೭ . ೧೯೨೨		בשת שמב	25 14:58:4	8 1993		21	
1277	126 VAL	и	51615	-42.26751	109.64790	Bl	44	ססססס. ס
1278	126 VAL		20939. در-	-42.12308	108.70456	B1	44	0.00000
1279			-34.59346	-41.07914	110.50496	Bl	4 4	0.00000
1280	126 VAL	CB		-40.82462			4.4	0.00000
1281	126 VAL	CG1	-36.22164	-39.41758	111.54094	B1	4.4	0.00000
1282	126 VAL	CG2	-37.04494	-41.002B1	109.82481	Bl	44	0.00000
1283	126 VAL		-33.57625	-41.13739	111.64961	Bl	4.4	0.00000
1284	126 VAL	. 0	-33.78819	-40.75401	112.79533	Bl	4 4 4 5	0.00000
1285	127 GLY	N	-32.41541	-41.66443	111.25324	B1	45	0.00000
1286	127 GLY	H		-42.01122		10 J	45	0.00000
1287	127 GLY	CX		-41.88201			45	0.00000
1288	127 GLY	C		-41.56889		BI	45 •	0.00000
1289	127 GLY	0	-29.03055	-41.21776 -41.65704			46	0.00000
1290 1291	128 GLU 128 GLU	Н	-29.01143 -30 48653	-42.13586			46	0.00000
1292	128 GLU	CA	-28.56256	-41.14297	109.73120		46	0.00000
1293	128 GLU	CB		-42.29481		Bl	46	0.00000
1294	128 GLU	CG	-26.29316	-42.25781	109.59860		46	0.00000
1295	128 GLU	CD		-43.28422			46	0.00000
1296	128 -GLU	- OE1-		-42.88073			46	0.00000
1297	128 GLU	OE2	-25.71010	-44.47773	109.00348		46	0.00000
1290	128 GLU	С		-39.98012			46	0.00000
1299	128 GLU	0		39.57375	108.35724		46	0.00000
1300	129 TYR	N		-39.42619			47	0.00000
1301	129 TYR	н	-26.68498	-39.83420	108.68820		47	0.00000
1302	129 TYR	CA.	-27.50019	-38.29454	107.46434	BI	47 47	0.00000
1303	129 TYR	CB	-26,63842	-37.15326		81 21	47	0.00000
1304	129 TYR	ÇG	-27.30857	-36.34222	110.34811		47	0.00000
1305	129 TYR	CDj	-26.67181	-36.21958 -35.42131	111.35148		47	0.00000
1306	129 TYR 129 TYR	CE1 CD2	-27.25566 -20.52827	-35.67832	108.83210		47	0.00000
1307 1308	129 TYR 129 TYR	CE2		-34.88063			. 47	0.00000
1309	129 TYR	CZ	-28.47327	-34.75726	111.08860		47	0.00000
1310	129 TYR	OH	-29.05005	-33.98459	112.07221		47	0.0000
1311	129 TYR	нн		-33.40059	111.67732		47	0.00000
1312	129 TYR	Ċ	-26.82531	-38.64384	106.15941	Вl	47	0.00000
1313	129 TYR	0	-25.66697	-39.04407	106.10492		47	0.00000
1314	130 ARG	N	-27.55686	-38.38162	105.08581		48	0.00000
1315	130 ARG	Н		-38.10451	105.16735		48 48	0.00000
1316	130 ARG	CA	•	-38.41254	103.80227		48	0.00000
1317	130 ARG	CB		-38.84699 -39.00526	101.34112	R1	48	0.00000
1318	130 ARG 130 ARG	CG	-27.21143 -28.23975	-39.41974	100.29808	Bl	48	0.00000
1319 1320	130 ARG 130 ARG	.NE		-39.38333	.98.95629		48	0.00000
1321	130 ARG	HE		-38.85034	98.82833		48	0.00000
1322	130 ARG	CZ		-39.96348	97.93202	Bl	48	0.00000
1323	130 ARG	NHI	-27.82365	-39.81313	96.69917		48	0.00000
1324	130 ARG			-40.23513	, 95.90996	Bl	48	0.00000
1325	130 ARG		-27.01064		96.53955	Bl	48	0.00000
1326	130 ARG	NH2		-40.6B730	98.14392		48 48	0.00000 0.00000
1327	130 ARG		-29.90446	-41.10492	97.39118		48	0.00000
1328	130 ARG			-40.81930	99.07957 103.45986		48	0.00000
1329	130 ARG	C		-37.06053 -36.11772	103.43988		48	0.00000
1330 1331	130 ARG	0		-36.99899	103.58668		49	0.00000
1331	131 ALA 131 ALA	и Н	-24 45060	-37.77908			49	0.00000
1333	131 ALA	CA	-24.28607	-35.84894	102.98902	ві	49	0.00000
1334	131 ALA	CB		-35.44271	103.80801	ы	49	0.00000
1335	131 ALA	c		-36.20633	101.58658		49	0.00000
1336	131 ALA	ō	-23.17892	-37.20532	101.36065		49	0.00000
1337	132 VAL	N	-24.28495		100.63154	Bl	50	0.00000 0.00000
:338	132 VAL	1;		-34.54695	100.85567	ום ים	50 50	0.00000
1339	132 VAL	CA		-35.79619	99.24192 98.36560	B 1	50 . 50	0.00000
1340	132 VAL	CB	-25.1257R	-35.11200	90.30300	۱ ت	50	

./DR1	MIN2.C	22	Tau Feb	25 14:58:4	8 1993	22	
1341	132 V	AL CG1	5.0250	5 -35.45687			
1342		AL CG2		8 - 35.48853			
1343		አኒ C		6 -35.53393			
1344		al o		B -36.19257	97.84114		
1345		HR N		6 -34.53567		-	
1346		HR H	-22.4403	6 -34.04272 7 -34.13851	98.91128		
1347		HR CA HR CB	-20.6623	9 -33.09962	97.76600		
1348 1349		HR OG1	-19.5851	-32.56706	97.34397		0.0000
1350	133 7		-19.6901	-32.10136	96.50551		
1351	133 T	_	-21.8173	-31.96947	98.12788		
1352	133 T		-19.9173	-33.59750	100.10846	B1 51	
1353	133 T	HR O	-20.48697		101.15881	B1 51	
1354	134 G		-18.60119	-33.44216	99.91793		
1355	134 G		-18.20352	-33.65725	99.02398		
1356	134 G		-17.75238		100.97647		
1357	134 G		-16.30452	2 -32.81284 3 -34.17073	100.43462		
1358	134 GI		-15.76445	-34.04410	99.60488		
1359 1360	134 GI	SU CD	-14.513// 13 54341	-34.96158			
1361	134 GI		-13.96021	-33.03519	98.99487	B1 52	0.0000
1362	134 GI		-18.18601	-31.51310	101.45728	B1 52	
1363	134 GI	_	-17.97884	.,-31.11916	102.59352	B1 52	
1364	135 LE		-18.87611	-30.80674	100.55812	B1 53	
1365	135 LE		-18.92762	-31.13340	99.61398	B1 53	
1366	135 LE	CU CA	-19.55727	-29.57029			
1367	135 LE		-20.29914	-29.06218	99.70959		
1368	135 LE		-20.12211	-27.57937	99.38843		
1369	135 LE		-20.93390	-26.68772	100.32666		
1370	135 LE		-18.63204	-27.23870	99.36436		
1371	135 LE		-20.53099	-29.72915 -28.93514	102.11322		
1372	135 LE		-21.29082		102.02891	E1 54	
1373 1374	136 GI 136 GI		-21 14062	-31.49656	101.30480	B1 54	0.00000
1375	136 GI		-22.25373	-31.10000	103.08935	B1 54	
1376	136 GI	•	-21.66227	-31.85099	104.26834	B1 54	
1377	136 GL		-22.17280	-31.82013	105.38053	B1 54	
1378	137 AR		-20.54192	-32.53644	104.01306	B1 55	
1379	137 AR	•	-20.17033	-32.57514	103.08296	B1 55 B1 55	
1380	137 AR	.*	-19.94165	-33.32437	105.09532	B1 55	
1381	137 AR		-18.67070	-34.02762 -35.07893	105.55510		
1382	137 AR		-18.13008	-36.15719	105.35023		
1383	137 AR		-19.10033	-37.06640	106.87657		
1384 1385	137 AR 137 AR		-17 93002	-36.71980	107.47204	B1 55	
1386	137 AR		-19.21896	-38.26612	107.07086	B1 55	0.00000
1387	137 AR		-18.67934	-39.07643	107.97489	B1 55	
1388	137 AR	G HR11	-19.03969	-39.99255	108.14749	B1 55	
1389	137 AR	G HH12	-17.88763	-38.75541	108.49972	B1 55	
1390	137 ÄR	G NH2	-20.29085	-38.64369	106.37214	B1 55	
1391	137 AR	G HH21	-20.72192	-39.53705	106.49834	B1 55	
1392	137 AR		-20.68879	-38.01490 -32.65026	105.70437		
1393 1394	137 AR		-17.70330	-33.13638	107.44172		
1395	138 PR		-19 03095	-31.51331	106.56277	B1 56	0.0000
1396	138 PR		-18.26903	-30.78028	105.55482	B1 26	
1397	138 PR		-10.92500	-30.86885	107.87839	B1 56	
1398	138 PR		-18.00966	-29.66775	107.61418	51 56	
1399	138 PR		-17.22921	-30.01427	106.35388	21 56	
1400	138 PR	•	-20.24853	-30.39451	108.45452	21 56	
1401	138 PR		-20.38873	-30.20105	109.652;3	31 56	
1402	139 AS		-21.22822	-30.19487	107.56513	B1 57	
1403	139 AS		-21.12955	-30.45258	100.00235	21 57	
1404	139 AS	P CA	-22.54445	-29.76374	100.04027	,	

./221	_KID:2 . CRO	•	בשה צפף	25 14:58:	48 1993		23	
1405	139 ASP	СВ	3 29483	-29.2279	4 106.80729	B1	57·	~0~0°0000
1406	139 ASP			-28.4415			57	0.00000
1.07	139 ASP		-24.33925	-27.2595	5 107.53698	81	57	0.00000
1408	139 ASP	OD2	-25.59590	-29.0093	7 107.24893	B1	57	0.00000
1409	139 ASP	С		-30.9065			57	0.00000
1410	139 ASP		-23.84428	-30.7994	109.79626		57	0.00000
1411	140 ALA		-23.18867	-32.06950	0 108.04975 5 107.13514		58 58	0.00000
1412	140 ALA			-32.09325 -33.29863			58	0.00000
1413 1414	140 ALA 140 ALA			-34.5128			58	0.00000
1415	140 ALA		-22.96994	-33.52593	110.04660		58	0.00000
1416	140 ALA			-33.62460			58	0.00000
1417	141 GLU		-21.63141	-33.53710	110.01537	Bl	59	0.00000
1418	141 GLU	H	-21.14260	-33.44422	109.14251	Bl	59	0.00000
1419	141 GLU	CA		-33.73045			59	0.00000
1420	141 GLU	СВ	-19.38545	-33.75474	110.92637		59 59	0.00000
1421	141 GLU	CG	-19.08157	-34.92155 -34.91070	109.97183		59	0.00000
1422	141 GLU 141 GLU	CD OE1	-17.65605	-34.91070	109.44077	B1	59	0.00000
1424	141 GLU	051	·· -16.99658	-33.87375	-109.50288	Bl	59	0.00000
1425	141 GLU	C	-21.20315	-32.72395	112.36696	Bl	59	0.00000
1426	141 GLU	ō	-21.35204	-33.05368	113.53869	Bl	59	0.00000
1427	142 TYR	N	-21.39109	31.47026	.111,93865	Bl	60	0.00000
1428	142 TYR	н		-31.22939			60	0.00000
1429	142 TYR	CA		-30.42572			60	0.00000
1430	142 TYR	CB	-22.17510	-29.18770	111.95478		60 60	0.00000
1431	142 TYR	CG		-27.86866 -27.21121			60	0.00000
1432	142 TYR 142 TYR	CD1 CE1	=20.91930	-27.21121			60	0.00000
1433	142 TYR	CD2		-27.27306			€0	0.00000
1435	142 TYR	CE2	-23.33600	-26.00211	113.72688	Bl	60	0.00000
1436	142 TYR	CZ	-22.10013	-25.34007	113.91274	B1	60	0.00000
1437	142 TYR	ОН	-22.06472	-24.08718	114.49226		60	0.00000
1438	142 TYR	нн	-22.95958	-23.75980			60 60	0.00000 0.00000
1439	142 TYR	Ç		-30.84932	113.52485		60	0.00000
1440	142 TYR	0	-23.33185	-30.86335 -31.24530	114.74427		61	0.00000
1441	143 TRR 143 TRP	n N	-24.16619	-31.27353	111.69669	Bì	61	0.00000
1443	143 TRP	CA	-25 46084	-31.65772	113.24427	Bl	61	0.00000
1444	143 TRP	CB	-26.46502	-31.82534	112.10045	Bl	61	0.00000
1445	143 TRP	CG	-26.82927	-30.51319	111.43167		61	0.00000
1446	143 TRP	CD2		-30.35383			61	0.00000
1447	143 TRP	CE2	-27.68725	-28.88192	110.01923	B1	61 61	0.00000
1448 1449	143 TRP.	CE3	-28.23171	-31.24445 -29.20408	111 84528	B1	61	0.00000
1450	143 TRP 143 TRP	CD1 NE1	-26.99373	-28.24161	111.01939	Bl	61	0.00000
1451	143 TRP	HE1	-26.86335	-27.27405	111.09738	Bl	61	0.00000
1452	143 TRP	CZZ	-28.41151	-28.41903	108.90296	81	61	0.00000
1453	143 TRP	CZ3	-28.94655	-30.73482	108.27096	Bl	61	0.00000
1454	143 TRP	CH2	-29.03488	-29.34388	108.03833	B1	61	0.00000 0.00000
1455	143 TRP	С	-25.40824	-32.93379	114.07770	BI	61 61	0.00000
1456	143 TRP	0 .	-26.13451	-33.11650	113.04993	BJ DT	62	0.00000
1457 1458	144 A6N	N	-24.46546 -23.94027	-33.80055	112 85783	B1	62	0.00000
1459	144 ASN 144 ASN	Н СА	-23.94047	-33.82303 -34.99080	114.49069	Bl	62	0.0000
1460	144 ASN	CB	-23.20850	-35.93308	113.73882	B1	62	0.00000
1451	144 ASN	CG	-23.89541	-36.68861	112.61740	Bl	62	0.00000
1462	144 ASN	ODl	-23.68155	-36.47075	111.43371	Bl	62	0.00000
1463	144 ASN	ND2	-24.72776	-37.64097	113.02365	Bl	62	0.00000
1464	144 ASN	HD21	-24.89338	-37.61926	113.99276	81	62 63	G.COOOO G.OOOOO
1465	144 ASN		-25.19584 -23.49875	-39.20406	114 P7591	B)	62 62	0.00000
1466	144 ASN	0	-23.49875	-34.03431	116.69934	91	62	0.00000
1458	144 ASN 145 SER	16	-23.43003	-33.46640	115.97353	31	63	0.00000
	270 000			,,,,,,,,	- :			

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1469	145 EE	R H	22.9783	0 -32.8042	В 115.21961 В1	65 **	
1470	145 SEF	R CA			1 117.25865 B1	63	0.00000
3473					4 117.01176 B1	63	0.00000
1472					0 118.24511 B1	63	0.00000
1473					5 118.91065 B1	63	0.00000
1474	145 SEF				4 118.06161 B1 7 119.13941 B1	63 63	0.00000
1475 1476	145 GLN				7 117.51252 B1	64	0.00000
1477	146 GLN				4 116.64774 B1	64	0.00000
1478	146 GLN				6 118.19522 Bl	64	0.00000
1479	146 GLN				1 117.23892 B1	64	0.00000
1480	146 GLN		-25.4585	5 -27.9694	5 117.72882 B1	64	0.00000
1481	146 GLN		-26.8909	6 -27.81462	2 117.25234 B1	64	0.00000
1482	146 GLN				9 118.00911 B1	64	0.00000
1483	146 GLN				3 115.94329 Bl	64	0.00000
1484 1485	146 GLN 146 GLN				9 115.33276 B1 7 115.56134 B1	64 64	0.00000
1486	146 GLN				118.57642 B1	64	0.00000
1487	146 GLN	ō			117.77294 B1	64	0.00000
1488	147LYS		26.36947	7-34-44426	119.82371 B1	65	0.00000
1489	147 LYS	н	-25.65249	-31.20699	120.47663 B1	65	0.00000
1490	147 LYS	CA			120.21964 B1	65	0.00000
1491	147 LYS	CB			121.37895 B1	65	0.00000
1492	147 LYS	CG			121.84203 B1	65 (5	0.00000
1493	147 LYS	CD CE			121.97244 B1 121.47444 B1	65	0.00000
1494 1495	147 LYS 147 LYS	NZ	•	-36.81034	•	65	0.00000
1496	147 LYS	HZ1		-37.42223		65	0.00000
1497	147 LYS	HZ2		-36.76558		65	0.00000
1498	147 LYS	HZ3			122.15073 B1	65	0.00000
1499	147 LYS	Ċ			119.09340 Bl	65	0.00000
1500	147 LYS	0	-29.30235	± 32.90174	118.83883 Bl	65	0.00000
1501	148 ASP	N		-34.00916		66	0.00000
1502	148 ASP	H		-34.14495		6 6 6 6	0.00000
1503 1504	148 ASP	CA CB		-35.00236 -36.37833		66	0.00000
1505	148 ASR	CG		-36.57930	119.21585 B1	66	0.00000
1506	148 ASP	OD1	-25.69280	-36.06325	119.05193 B1	66	0.00000
1507	148 ASE	OD2		-37.20742	120.23524 B1	66	0.00000
1508	148 ASP	C .		-34.63043		66	0.00000
1509	148 ASP	0	-30.17136	-35.14808	117.04951 B1	66	0.00000 0.00000
1510 1511	149 LEU 149 LEU	n H		-33.61840 -33.30949	115.93941 B1 115.73440 B1	67 67	0.00000
1512	149 LEU	CA	-27.36365	-32.84836	115.45008 B1	67	0.00000
1513	149 LEŲ	СВ	-29.45399	-31.74267	114.57360 B1	67	0.00000
1514	149 LEU	CG	-30.29432	-30.58223	114.02475 B1	67	0.00000
1515	149 LEU	CD1	-30.85820	-30.92475	112.65290 B1	67	0.0000
1516	149 <u>L</u> EU	CD2			115.02072 B1	67	0.00000
1517	149 LEU	Č			114.80952 B1	67	0.00000 0.00000
1518	149 LEU	0	-32.36472	-33.37859	114.83807 Bl 114.32363 Bl	67 68	0.00000
1519 1520	150 LEU 150 LEU	H jv	-29 75579	-34 94139	114.11621 B1	68	0.00000
1521	150 LEU	CA	-31.59782	-36.02822	114.12850 B1	68	0.00000
1522	150 LEÚ	CB	-30.74740	-37.29867	114.15286 B1	68	0.00000
1523	150 LEU	CG	-29.89363	-37.44772	112.89569 B1	68	0.00000
1524	150 LEU	CD1	-20.64060	-38.26626	113.18796 B1	68	0.00000
1525	150 LEU	CD2	-30.71709	-38.01712	111.73915 B1	68	0.00000
	150 LEU	C			115.10785 B1	65 -63	0.00000 0.00000
1527 1528	150 LEU 151 GLU	0 N			114.70350 B1 116.39428 B1	63	0.00000
	151 GLU	H	-31.52943	-36.57284	116.72119 B1	6.5	0.00000
1530	151 GLU	CA	-33.61295	-36.59512	117.30950 EL	69	0.00000
1331	151 GLU	CB	-33.19489	-36.98331	118.72928 Bl	65	0.00000
1532	151 GLU	CG	-32.69081	-38.41906	118.86324 91	6 9	0.00000

. אנו	TEXTHS . CKI	o	Thu Fo	ය 25 14:58	:48 1993	25	
153		J CD	-31,192	87 -38.444	64 118.70497 E	1 69	0.00000
135	s 151 GLU	J OE1	-30.713	20 -30.331	29 117.58148 B	1 69	0.00000
153		OE2	-30.515	22 -38.564	62 119.72459 B	1 69	0.00000
153					68 117.43857 B		0.00000
153			-35.757	90 -35.573	13 117.58647 B	1 69	0.00000
1538					80 117.35586 B		0.00000
1539 1540					56 117.18602 B		0.00000
1543					52 117.40695 B 46 117.32018 B		0.00000
1542					00 118.60094 B		0.00000
1543					24 118.49387 B		0.00000
1544	152 GLN	OE1 -	-32.130	68 -29.270	19 118.36008 B	70	0.00000
1545					52 118.54711 B		0.00000
1546					11 118.65663 B		0.00000
1547 1548					58 118.46911 B		0.00000
1549					16 116.27182 BI N1 116.43607 BI		0.00000
1550					69 115.08552 BI		0.00000 0.00000
1551	153 ARG				9 114.99299 BI		0.00000
1552		· · CA · · · -	36.2585	3 -33:4212	9 119.95947 B1	71	0.00000
1553	153 ARG	CB -	35.4632	2 -33.4370	7 112.66564 BI	71	0.00000
1554	153 ARG	ce -	34.8428	0 -32.0779	1 112.35054 B1	71	0.00000
1555	153 ARG				0.111.17839 B1	71	0.00000
1556	153 ARG				6 110.55166 B1	71	0.0000
1557	153 ARG	HE -			O 110.98860 B1	71	0.00000
1558 1559	153 APG				2 109.31716 B1	71	0.00000
1560	153 ARG 153 ARG				6 108.54980 B1 2 107.55422 B1	71 71	0.00000 0.00000
1561	153 ARG				2 107.33422 B1 4 108.90136 B1	71	0.00000
1562	153 ARG				1 108.83735 B1	71	0.00000
1563	153 ARG				2 107.04355 B1	71	0.00000
1564	153 ARG	нн22 -:	32.4557	1 -32.9735	9 109.42307 B1	71	0.0000
1565	153 ARG				B 114.06926 B1	71	0.00000
1566	153 ARG				B 113.96948 Bl	71	0.00000
1567 1568	154 ARG				114.34589 B1	72	0.00000
1569	154 ARG 154 ARG				5 114.34685 B1	72	0.00000
1570	154 ARG				3 114.59374 B1 3 115.12926 B1	72 72	0.00000 0.00000
1571	154 ARG				115.22908 B1	72	0.00000
1572	154 ARG				115.69023 B1	72	0.00000
1573	154 ARG				115.80184 B1	72	0.00000
1574	154 ARG		7.59626	-42.01330	115.09038 Bl	72	0.00000
1575	154 ARG		**		116.80504 B1	72	0.00000
1576 1577	154 ARG 154 ARG				116.88687 Bl	72	0.00000
1578	154 ARG	NU11 -3	9 17405	-44.4/83/ -/3 06000	117.61587 B1 116.20728 B1	72 72	0.00000
1579	154 ARG				117.71513 B1	72	0.00000
1580	154 ARG				118.48012 B1	72	0.00000
1581	154 ARG	HH22 -3	5.18100	-41.66590	117.63709 B1	72	0.00000
1582	154 ARG				115.51516 B1	72	0.00000
2583					115.28503 B1	72	0.00000
1584					116.55472 B1	73	0.00000
1585 1586					116.77047 B1	73 73	0.00000
					117.36371 B1 118.48265 B1	73 73	0.00000 0.00000
		CS 4 (0.15723	-34.36970	118.46265 B1 119.44775 B1	73	0.00000
1589					120.46128 31	73	0.00000
1590					121.28237 B1	. 73	0.00000
1591	155 ARG	HE - 41	63763	-33.18145	121.09517 81	73	0.00000
					122.22743 81	73	0.00000
					122.94081 B1	73	0.00000
					123.65592 El	73 73	0.00000 0.00000
					122.77039 B1 122.45304 B1	7.7	0.00000
	1			J 4 7 5 0 9		• •	

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1661	162 ARC	311	50.49694	-40.8845	114.11092	Bl	٤O	σ.00000
1662	162 ARG				114.80601		80	0.00000
ر ت	162 ARG				114.46730		80	0.00000
1664	162 ARG	: нн13	-50.11979	-44.31403	114.66150	B1	80	0.00000
1665			2 -51.06184	-43.06285	113.96028	Bl	80	0.00000
1666		NH2	-48.19839	-42.76163	115.38323	Bl	80	0.00000
1667		нн21	-48.00134	-43.72771	115.55650	B)	80 80	0.00000
1668			2 -47.52579	75 02006	115.64161 113.42487	DI TG	80	0.00000
1669			-49.75514 -50.02003	-35.92900	112.81658	B1	80	0.00000
1670 1671	163 HIS		-30.02093	-35.33752	114.60088	Bl	81	0.00000
1672	163 HIS		-48 68252	-35.28035	115.03933	B1	81	0.00000
1673	163 HIS		-50.76431	-34.73357	115.23467	Bl	81	0.00000
1674	163 HIS		-50.42874	-34.29045	116.66134	Bl	81	0.00000
1675	163 HIS		-50.41026	-35.50649	117.55950	B1	81	0.0000
1676	163 HIS	ND1	-51.51644	-36.01842	118.12240	Bl	81	0.00000
1677	163 HIS		-52.42830	-35.67431	118.03242	B1	81	0.00000
1678	· 163 HIS	CD2	-49.31404	-36.28837	117.93420	B1	81	0.00000 0.00000
1679	163 HIS	NE2	-49.78030	-37.28089	118.73130	B1	81 81	0.00000
1680	163-HI6 163 HI6		-51.13923	-37 505RA	-118.84798 114.44466	Bl	81	0.00000
1681 1682	163 HIS	С 0	-51.37160	÷33.55956	114.18659	B1	81	0.00000
1683	164 ASN	Ŋ		-32.67616			82	0.0000
1684	164 ASN	н	-49.51981	-32.74462	114.21247		82	0.00000
1685	164 ASN	CA	-51.04300	-31.58239	113.20331	91	82	0.00000
1686	164 ASN	CB	-49.96583	-30.54361	112.89196	Bl	82	0.00000
1687	164 ASN	CG.	-49.91907		114.02377		82	0.00000
1688	164 ASN	OD1.					82	0.00000
1689	164 ASN	ND2	-50.80473	-28.55321	113.91309	Bl	82	0.00000
1690	164 ASN	HD21	-51,44322	-28.49843	113.14586	Bl	52	0.00000
1691	164 ASN		-50.84210				82	0.00000 0.00000
1692	164 ASN	C		-32.02333			82 82	0.00000
1693	164 ASN	0	-52.73418 -51.12326	-31.48002	111.51085		83	0.00000
1694 1695	165 TYR 165 TYR	N	-50.28201		111.62194		83	0.00000
1696	165 TYR	Н СА	-51.81304		110.10401		83	0.00000
1697	165 TYR	CB	-50.92781		109.25048		83	0.00000
1698	165 TYR	CG	-51.39689	-34.50284	107.80317	Bl	83	0.00000
1699	165 TYR	CD1	-52.18946	-33.44669	107.29137		83	0.00000
1700	165 TYR	CEl	-52.63312	-33.47413	105.96126		83	0.00000 0.00000
1701	165 TYR	CD2	-51.04289	-35.58963	106.96903	22 D.T	83 83	0.00000
1702 1703	165 TYR 165 TYR	CE2 CZ	-51.48846 -52.28557	-35.61344			83	0.00000
1703	165 TYR		-52.75931	-34.53152	103.84155	B1	83	0.00000
1705	165 TYR	нн	-52.13515			Bl	83	0.00000
1706	165 TYR	Ç	-53.16114	-34.17050	110.39688	B1	83	0.00000
1707	165 TYR	0	-54.17243	-33.75040	109.85354	Bl	83	0.00000
1708	166 GLY	N	-53.16827	-35.13255	111.32677	B1	84	0.00000
1709	166 GLY	H	-52.30920	-35.44284	111.74443	B1	84	0.00000
1710	166 GTX	CA	-54.44388	-35.75931	111.69489	B1	84	0.00000 0.00000
1711	166 GLY	C	-55.55421	-34.78683	112.08191	81 81	84 84	0.00000
1712	166 GLY	0	-56.70058	-34.86763	111.64884	B1	85	0.00000
1713 1714	167 VAL 167 VAL	N H	-22.10423	-33.01049	113.25020	Bl	85	0.00000
1717	167 VAL	CA	-56.14083	-32.77588	113.26808	B1	8.5	0.00000
1716	167 VAL	CB	-55.54422	-31.87200	114.36486	51	85	0.00000
1717	167 VAL	CG1	-56.46931	-30.71890	114.75691	B1	8.5	0.00000
1718	167 VAL	CG2	-55,22649	-32.70528	115.60849	Bl	85	0.00000
1719	167 VAL	С	-56.62003	-31.96437	112.06515	B1	85 05	0.00000 0.00000
1720	167 VAL	0	-57.80658	-31.70971	111.87365	EJ ET	85 86	0.00000
1721	168 GLY	N	-55.65605	-31.61592	111.20320	E 3	86	0.00000
1722	168 GLY	H	-34.6862/ -56.04316	-31.01344 -30.44480	111.38362 109.94972	Bl	86	0.00000
1723 1724	158 GLY 158 GLY	CA C	-50.04213 -57 02338	-31.77629	109.11607	B 1	86	0.00000
-124	-20 GF:	•	57.02550				_	

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1725	168 GL	. 0	58.02649	-31,28044	108.61617	Bl	86	0.00000
1726				-33.07169			87	0.00000
1717					109.43169		87	0.00000
1728		CA			108.30469		87	0.00000
1729					108.28846		87	0.00000
1730					107.43481		87	0.00000
1731					107.65905		87 87	0.00000
1732				-30.01342	108.02205	B1	87	0.00000
1733 1734			-50.40540	-34 N5273	108.86565	B1	87	0.00000
1735					108.14126		87	0.00000
1736	170 SER		-59.06996	-33.99305	110.19884	Bl	88	0.00000
1737	170 SER	-	-58.23138	-34.05998	110.74831	Bl	88	0.00000
1738	170 SER		-60.38255	-33.85094	110.83391	Bl	88	0.00000
1739	170 SER	CB			112.35798		88	0.00000
1740	170 SER				113.03659		88	0.00000
1741	170 SER		-61.30877	-34.00943	113.98374	B1	8 8 8 8	0.00000
1742 1743	170 SER 170 SER				110.37646		88	0.00000
1744	171 SER		-62.31497 -60.40231	-32.69130	-110.41676	B1	89	0.00000
1745	171 PHE	н	-59.54601	-31.41832	110.75393	Bl	89	0.00000
1746	171 PHE	CA	-61.19539	-30.24631	109.95663	Bl	89	0.00000
1747	171 PHE	CB	-60.30793		110.10880		89	0.00000
1748	171 PHE	CG	-59.94208	-28.68147	111.54294	Bl	8 9	0.00000
1749	171 PHE	CD1	-58.59291		111.85413		89	0.00000
1750	171 PHE	CD2	-60.93098		112.55491		89	0.00000
1751	171 PHE	CEI	-58.23032		113.17656		89	0.00000
1752	171 PHE	CE2	-60.56845		113.87858		89	0.00000
1753	171 PHE	CZ	-59.21901		114.18527		89	0.00000
1754	171 PHE	С	-61.62802		108.49502		89 89	0.00000
1755 1756	171 PHE 172 THR	О. И.	-62.68697 -60.72520		107.69903		90	0.00000
1757	172 THR	И, Н	-59.91792		108.09636		90	0.00000
1758	172 THR	CA	-60.84308		106.25599		90	0.00000
1759	172 THR	CB	-59.41710		105.70165	B1	90	0.00000
1760	172 THŔ	0G1	-59.35292	-30.21143	104.36207		90	0.00000
1761	172 THR	HG1	-60.15430		103.89285	B1	90	0.00000
1762	172 THR	CG2	-58.90262		105.74825		90 90	0.00000
1763 1764	172 THR	Ċ	-61.71208 -61.76078		105.47987		90	0.00000
1765	172 THR 173 VAL	О	-62.32957		106.19315		91	0.00000
1766	173 VAL	н	-62.28706		107.19363		91	0.00000
1767	173 VAL	CA	-62.87984		105.53718		91	0.00000
1768	173 VAL	· CB	-63.87967	-34.51911	106.47899		91	0.00000
1769	173 VAL	CG1	-64.50006	-35.77324	105.85034	Bl	91	0.00000
1770	173 VAL	CG2	-63.20452	-34.90248	107.79466	B1	91	0.00000
1771	173 VAL	С	-63.51710	-33.59419	104.16716	B 1	91	0.00000
1772	173 VAL	0	-63.18750	-34.24452	103.18126	B1	91 92	0.00000
1773 1774	174 GLN	N	-64.41211 - -64.59818 -	-32.59570	104.14/11	81 B1	92	0.00000
1775	174 GLN 174 GLN	Н СА	-65.14373 ·	-32.11269	103.00138	B1	92	0.00000
1776	174 GLN	C3	-65.76132 ·	-30.80951	103.22574	B1	92	0.00000
1777	174 GLN	CG	-66.77986 -	-30.26016	102.21318	Bl	92	0.00000
1778	174 GLN	CD	-66.09397		101.04012	B1	92	0.00000
1779	174 GLN	OE1	-65.45924		101.15380	Bl	92	0.00000
1780	174 GLN	NE2	-66.26402 -		99.87453		92	0.00000
1781	174 GLN	HE21	-66.63226		99.82928		92	0.00000
1782	174 GLN		-65.97833 -		59.02730		92 62	0.00000 0.00000
1783	174 GLN	С	-64.36067 -		101.63553		92 92	0.00000
1784	174 GLN	0	-64.88582 - -63.09233 -		100.57991		93	0.00000
1785 1786	175 ARG 175 ARG	И Н	-62.69754 -		102.62225	B1	93	0.00000
1787	175 ARG	r CA	-62.31707 -		100.50018	B 1	93	0.00000
1788	175 ARG	CB	-61.85817 -		100.16958		93	0.00000
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1789	175 ARG		-61 2/30	5 -30.26914	90.77310 B1	93	0.00000
1790				7 -28.04209		93	0.00000
1791	175 ARG			7 -28.87216		93	0.00000
1792	175 ARG			6 -29.76910		93	0.00000
1793	175 ARG			5 -27.74924		93	0.00000
1794	175 ARG			3 -27.83579		93	0.00000
1795	175 ARG		1 -58.1941	4 -27.02728	95.28229 B1	93	0.00000
1796	175 ARG	HH12	2 -58.22480	-28.73459	95.46034 Bl	93	0.00000
1797	175 ARG	NH2	-59.92427	7 -26.55682	97.09593 B1	93	0.00000
1798	175 ARG			3 -25.70796		93	0.00000
1799	175 ARG	HH22	2 -60.60313	-26.50975	97.82696 B1	93	0.00000
1800	175 ARG	C			100.46150 B1	93	0.00000
1801	175 ARG	0		-33.52738	99.49809 Bl	93	0.00000
1802	176 ARG	N		3 -32.85103		94	0.00000
1803	176 ARG	н		32.31340		94	0.00000
1804	176 ARG	CA		-33.80802	101.41499 Bl	94	0.00000
1805	176 ARG	CB		-33.62483	102.49829 B1	94	0.00000
1806	176 ARG	CG		-32.37173	102.34974 B1	94	0.00000
1807	176 ARG	CD		-32.28237	103.45854 Bl	94	0.00000
1808	176 ARG	NE		-30.94018	104.04511 B1	94	0.00000
1609	176 ARG	HΣ		-30.25818	103.53710 B1	94 94	0.00000
1810	176 ARG	CZ		-30.63659	105.19759 B1	94	0.00000
1811	176 ARG	NHI		:-29.38340	105.64628 B1	94	0.00000
1812	176 ARG	HH11		-29.10955 -28.69008	106.50132 B1 105.12278 B1	94	0.00000
1813 1814	176 ARG 176 ARG	нн12 Nн2		-31.57276	105.89297 B1	94	0.00000
1815	176 ARG	HH21		-31.37726	106.75753 B1	94	0.00000
1816	176 ARG			-32.51329	105.54660 B1	94	0.00000
1817	176 ARG	C		-35.27348	101.42902 B1	94	0.00000
1818	176 ARG	Ö		-36.15316	101.06767 B1	94	0.00000
1819	177 VAL	ห		-35.53390	101.85272 B1	95	0.0000
1820	177 VAL	н		-34.82991	102.18430 B1	95	0.00000
1621	177 VAL	CA	-61.32863	-36.91608	101.70299 B1	95	0.00000
1822	177 VAL	CB .	-61.79519	-37.47193	103.06780 B1	95	0.00000
1823	177 VAL	CG1	-61.84113	-39.00335	103.05589 B1	95	0.00000
1824	177 VAL	CG2	-60.88761	-37.00753	104.21299 Bl	95	0.00000
1825	177 VAL	С		-37.02458	100.63553 B1	95	0.0000
1826	177 VAI	0	-63.44404	-37.67295	100.79118 B1	95	0.0000
1827	178 XIS	N		-36.33028	99.52296 B1	96	0.00000
1828	178 HIS	н	•	-35.77698	99.39184 B1	96	0.00000
1829	178 HIS	CA		-36.32813	98.41659 B1	96	0.00000
1830	178 HIS	CB		-35.10064	98.57314 B1	96	0.00000
1831	178 HIS	CG		-35.21481	97.74669 B1	96	0.00000
1832		-ND1		-35.62002	96.46944 B1	96 96	0.00000
1833	178 HIS	HD1		-35.91855	95.96816 B1 98.14872 B1	96	0.00000
1834	178 HIS	CD2		-34.92566	98.14872 B1 97.08780 B1	96	0.00000
1835	178 HIS	NE2		-35.16505	96.05191 B1	96	0.00000
1836 1837	178 HIS 178 HIS	CE1 C		-35.59364 -36.29235	97.09131 B1	96	0.00000
1838	178 HIS		-61.22615		97.07091 B1	96	0.00000
1839	178 HIS		-62.87363		96.08759 Bl	96	0.00000
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FIG. 30

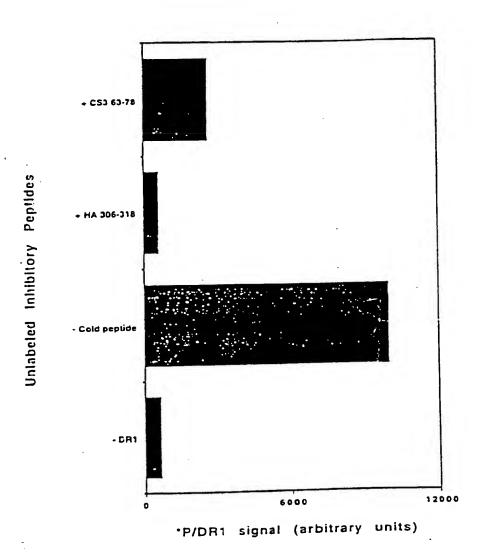
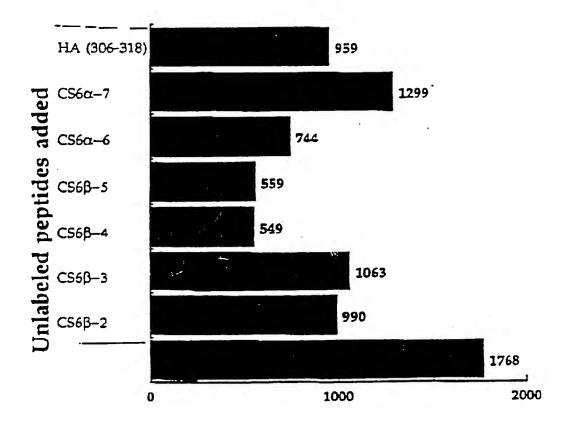


FIG. 31

Inhibition of 125 I HA (306-318)/DRI by unlabeled CSG of and B pertides



*HA/DR1 compact dimer signal (densitometric units)

FIG. 32

31/31

INTERNATIONAL SEARCH REPORT

International application No.
PCT/US94/05697

A. CLASSIFICATION OF SUBJECT MATTER IPC(5) : A61K 39/00, 39/02, 39/12, 37/02, 35/14 US CL : 424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75 According to International Patent Classification (IPC) or to both national classification and IPC B. FIELDS SEARCHED Minimum documentation searched (classification system followed by classification symbols) U.S. : 424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75 Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched Electronic data base consulted during the international search (name of data base and, where practicable, search terms used)						
C. DOCUMENTS CONSIDERED TO BE RELEVANT						
Category*	Citation of document, with indication, where ap	vant passages	Relevant to claim No.			
×	The Journal of Immunology, Volusies and 15 April 1993, Nauss et al Peptides in a Structural Homology MHC ", page 41A, Abstract 221,	1, 3-20				
Х Y	Nature, Volume 358, issued 27 A "Predominant Naturally Processed DR1 are derived from MHC-re Heterogenous in Size", pages 764-2, and Table 3.	l Peptides Bou lated Molecul	nd to HLA- e and are	12 1, 3-7		
X Furth	ner documents are listed in the continuation of Box C	. See pate	nt family annex.			
Special categories of cited documents: 'A' document defining the general state of the art which is not considered to be of particular relevance 'E' earlier document published on or after the international filing date document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified) 'O' document referring to an oral disclosure, use, exhibition or other means 'P' document published prior to the international filing date but later than the priority date claimed		"Y" later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention "X" document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step when the document is taken alone "Y" document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art document member of the same patent family				
Date of the actual completion of the international search O1 SEPTEMBER 1994 Date of mailing of the international search report 1 3 SEP 1994.						
Name and mailing address f the ISA/US Commissioner of Patents and Trademarks Box PCT Washington, D.C. 20231		Authorized fficer H. Sidberry	W. Ku;	za for		

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

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C (Continua	ation). DOCUMENTS CONSIDERED TO BE RELEVANT		
Category*	Citation of document, with indication, where appropriate, of the relevan	Relevant to claim No	
Y	The Journal of Immunology, Volume 150, No. 2, issued January 1993, Boehncke et al., "The Importance of Dor Negative Effects of Amino Acid Side Chain Substitution Peptide-MHC Molecule Interactions and T Cell Recognipages 331-341, see Abstract, on page 331.	8-11	
x	The EMBO Journal, Volume 9, No. 6, issued 1990, Jaral., "Peptide binding to HLA-DR1: a Peptide with most substituted to alanine retains MHC binding", pages 1797 page 1798, page 1800, figure 4, and page 1801, figure 2	512	
Ÿ	Nature, Volume 332, issued 28 April 1988, Brown et al hypothetical model of the foreign antigen binding site of histocompatibility molecules", pages 845-850, see pages	Class II	1, 3, 4
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Form PCT/ISA/210 (continuation of second sheet)(July 1992)*

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

Box 1 Observations where certain claims were f und unsearchable (Continuation of item 1 f first sheet)
This international report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:
Claims Nos.: 2 because they relate to subject matter not required to be searched by this Authority, namely:
Claim 2 is directed to a computerized model which encompasses scientific theory and computer programs to the extent that the International Searching Authority is not equipped to search prior art concerning such programs. Accordingly claim 2 is withdrawn from search under PCT Rule 39 and PCT Article 17(2)(a)(i).
Claims Nos.: because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:
Claims Nos.: because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).
Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)
This International Searching Authority found multiple inventions in this international application, as follows:
1. As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.
2. As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.
3. As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:
4. No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:
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Remark on Protest The additional search fees were accompanied by the applicant's protest.
No protest accompanied the payment f additional search fees.

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